Final Report Screening Site Inspection Laclede Coal Gas St. Louis, Missouri EPA ID# MOD981715980 TDD #F-07-9008-020 PAN #FM00579SA Site #Y33 Project #002 Prepared by E & E/FIT for the Region VII EPA RPO Project Manager: Keith Brown Superfund Contact: Greg Reesor Date: October 29, 1991

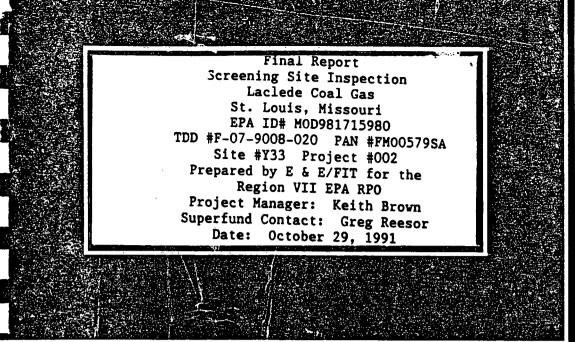
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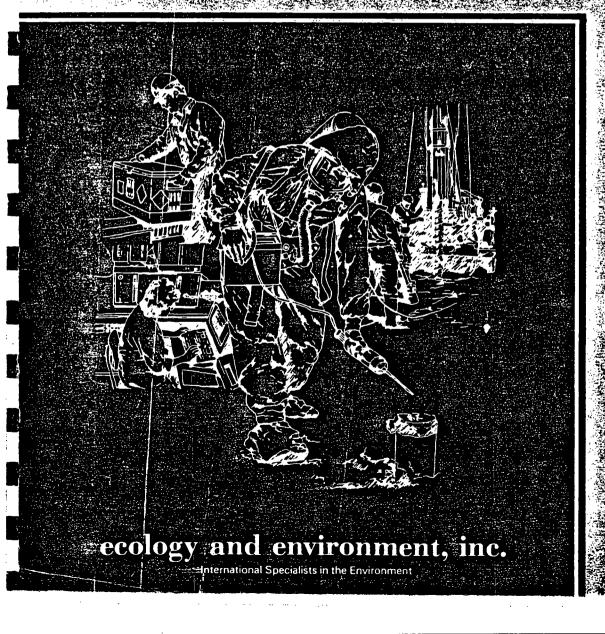
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SECTION 1: INTRODUCTION

The Ecology and Environment, Inc., Field Investigation Team
(E & E/FIT) was tasked by the Region VII U.S. Environmental Protection
Agency (EPA), through Technical Directive Document (TDD) #F-07-9008-020
(Appendix A) and EPA CERCLIS ID #MOD981715980, to conduct a Screening
Site Inspection (SSI) of the Laclede Coal Gas site located in St. Louis,
Missouri, foot of Mullanphy Street. The objectives of this SSI were to
determine the approximate areal and vertical extent of subsurface
contamination; to determine whether contamination is present in the 0 to
2 foot interval; to determine whether contaminants have migrated via
surface water runoff, or via the ground water/surface water interface;
and to determine whether any coal tar wastes present have leached into
the shallow ground water.

The E & E/FIT conducted a site reconnaissance on November 20, 1990, to assess current site conditions and facilitate preparation of the SSI work plan. Participating FIT members included Keith Brown, Bob Overfelt, and Chris Villiams. During this site reconnaissance, the FIT observed seepage from the foundation and piping system of an abandoned pump house that was formerly owned by the Mound Street Power Plant. Because this pump house is located on the east side of the flood control levee, this leachate was observed to be seeping directly into the Mississippi River.

During the SSI, the FIT collected subsurface soil, surface soil, surface water, ground water, sediment, and product samples. Initially, 24 locations were drilled with solid stem augers. Ninety-two samples were collected from these locations and screened on site in the mobile lab. Four of the sample locations determined by FASP analyses to be the most contaminated, as well as one background sample location, were resampled with hollow stem augers. Thirteen subsurface soil samples were collected from these 5 locations. Twelve surface soil samples were collected at 12 locations. Six surface water samples were collected at 4 locations. Six ground water samples were collected at 5 locations. Four sediment samples were collected at 3 locations. Two IDW samples were collected from 17 drums, and 1 oil sample was collected from an 80,000-gallon tank located on site. Additionally, two field/trip blanks, one auger/split spoon rinsate, and one Geoprobe pipe rinsate sample were

collected.

This report describes present site condition, dicusses site history and coal gas operations, the physical and cultural setting of the site, and presents and discusses the sample results of the FIT SSI. EPA Site Inspection Form 2070-13 was completed and is included as Appendix B.

SECTION 2: SITE LOCATION AND DESCRIPTION

2.1 SITE LOCATION

The Laclede Coal Gas site is located in St. Louis County, St. Louis, Missouri, approximately one mile north of the St. Louis Arch, along the Mississippi River at the foot of Mullanphy Street (Figure 2-1). The legal description of the site is city block 234-Tract #25, St. Louis Plan. The geographic coordinates of the site are 90° 11′ 00° west longitude, and 38° 38′ 20° north latitude. The site is situated in an industrialized area adjacent to the Mississippi River.

2.1 SITE DESCRIPTION

The site is a former manufactured gas plant (FMGP) (Figure 2-2). The dimensions of the site are approximately 600 feet by 500 feet. Currently, the site is owned and operated by Petroleum Fuels & Terminal Company (PF & T), a subsidiary of Apex Oil Company. PF & T operates a privately owned petroleum tank farm on site (Figure 2-3). The former Mound Street Power Plant building is owned by McKinley Iron and is located adjacent (northeast) to PF & T. The bulding was in the process of being raised during the SSI (E & E/FIT 1990).

The site is not secured and access to the ground is relatively unrestricted. There are locks on most of the doors of the site buildings and a fence surrounds the petroleum storage tanks. Also, there is a watchdog which lives within the fenced petroleum tank farm; no other security exists (E & E/FIT 1990).

The E & E/FIT conducted a site reconnaissance on November 20, 1990, to facilitate preparation of the work plan. The FIT observed seepage from the foundation and piping system of an abandoned pump house that we formerly owned by the Mound Street Power Plant. The pipes, which originate from the former plant, have been plugged with concrete, but seepage was still leaching through the concrete. Because this pump house is located on the east side of the flood control levee, this leachate was observed to be seeping directly into the Mississippi River.

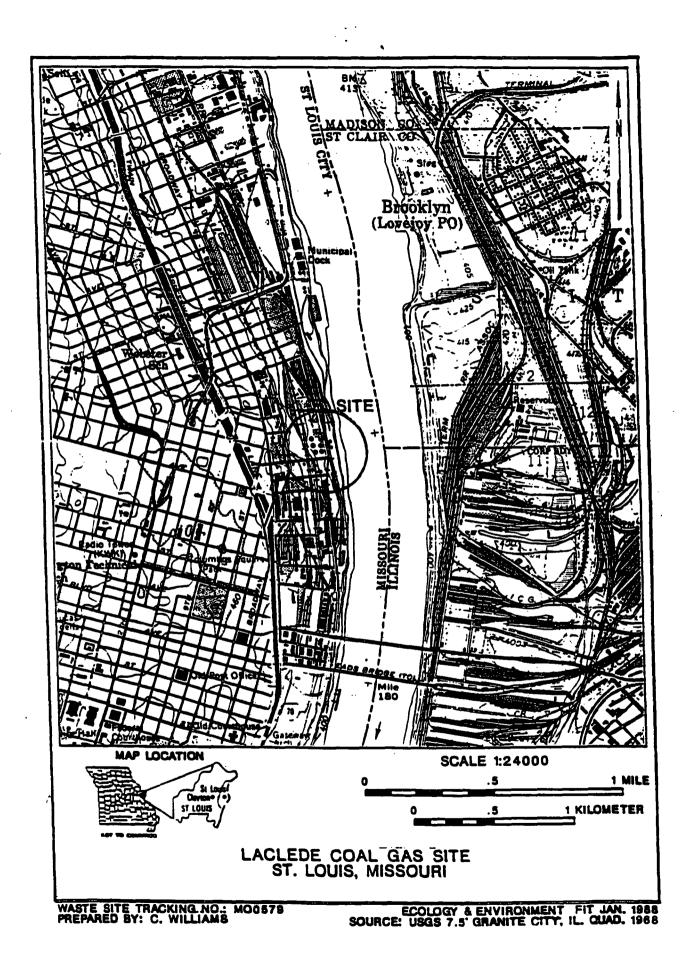
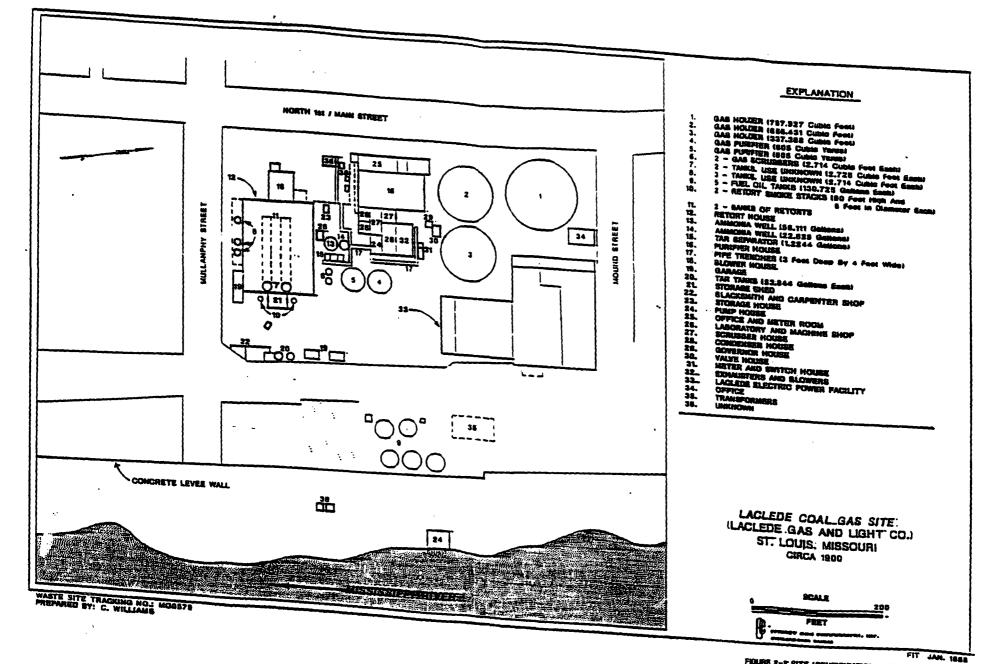
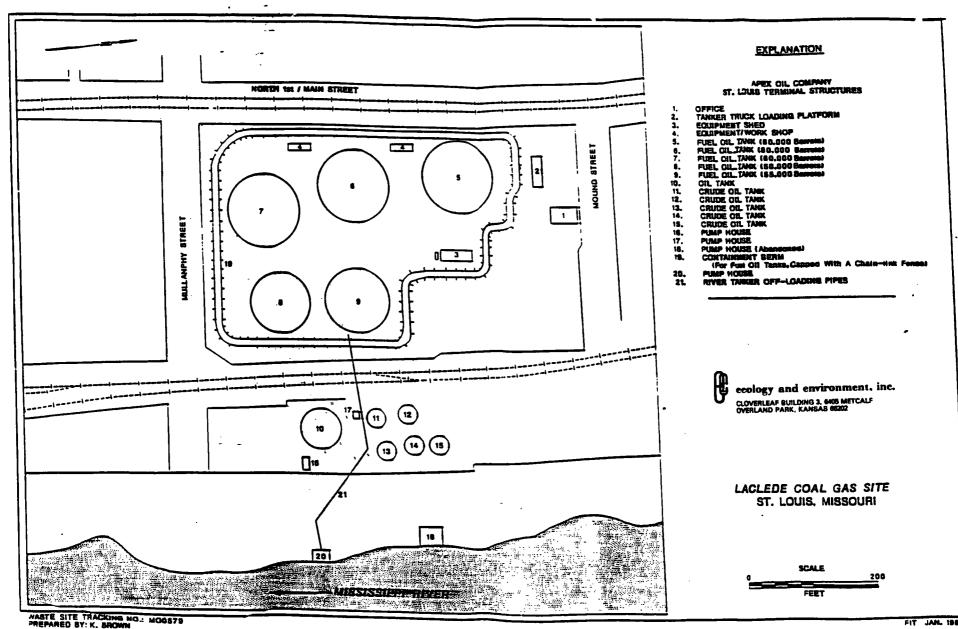


FIGURE 2-1: SITE LOCATION





2.3 SITE CONTACTS

Randel H. Lewis Terminal Manager Petroleum, Fuel & Terminal Company Foot of Mullanphy Street St. Louis, MO 63105 (314) 621-0522

Dick Bloch
Assistant Terminal Manager
Petroleum, Fuel & Terminal Company
Foot of Mullanphy Street
St. Louis, MO 63105
(314) 621-0522

Bob Welch Terminal Property Manager Petroleum, Fuel & Terminal Company Foot of Mullanphy Street St. Louis, MO 63:05 (314) 621-0522

Greg Reesor Missouri Coordinator U.S. Environmental Protection Agency Region VII 726 Minnesota Avenue Kansas City, KS (913) 551-7695

See Appendix C for a breakdown of samples by property owners and for addresses and phone number of other interested parties.

SETTON 3: SITE HISTORY

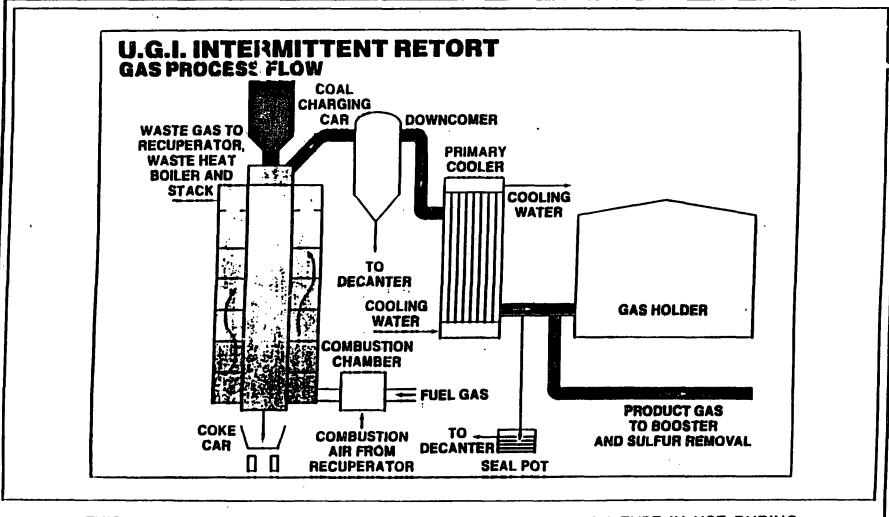
3.1 GENERAL SITE HISTORY

The earliest property records available indicate that this parcel of land was used by the Mound Street Warehouse Corporation, which sold the land and buildings to the Laclede Gas and Light Company on February 8, 1888 (E & E/FIT 1988). The Laclede Gas and Light Company constructed a large coal gasification plant on the property (Figure 2-2). According to the Brown's Directory of Manufactured Gas Plants, on March 23, 1940, the name of the facility was changed to the Laclede Power and Light Corporation (aka Laclede Electric), and the Laclede Gas Light Company (aka Laclede Gas). This suggests that the operations were separated, at least financially (Andrew 1890-1940).

Some time between 1940 and 1945, Phoenix Light, Heat and Power was involved in the Laclede operations; however, the extent of this involvement could not be determined during the background information search conducted for the work plan. On March 23, 1945, the entire facility was sold to Union Electric Company. According to Union Electric representatives. Union Electric Company never manufactured gas at this site (E & E/FIT 1988). Therefore, 1945 is the approximate closure date of the coal gasification works. Union Electric continued to use the electric power facility until 1973, though the former coal gas works was sold to the Apex Oil Company in 1969. Between 1969 and 1972, Apex Oil dismantled the old coal gas plant and constructed a petroleum tank farm on the site (Figure 2-3). This Apex facility stored various petroleum fuels until the mid 1980s when it became one of two Apex Oil asphalt product terminals in St. Louis. Currently, the terminal stores and distributes asphalt and #6 fuel oil (E & E/FIT 1988). Figure 2-2 depicts the Laclede Gas and Light Company as it appeared in the early 1900s; Figure 2-3 illustrates the current layout of the property.

3.2 FORMER SITE OPERATIONS

The major gas manufacturing process used at the Laclede FMGP was the UGI intermittent retort process (Environmental Research 1984)). This method produced gas through coal carbonization (Figure 3-1). During this process, coal is heated in the retort and the resulting coal gas is



THIS MANUFACTURED GAS MACHINE WAS THE MOST COMMON TYPE IN USE DURING THE PEROID OF OPERATION FOR THE LACLEDE FACILITY.

FIGURE 3-1

removed through its top. The gas is run through a condenser and a scrubber before it is moved into the gas holder. Wastes are produced in the condenser and scrubber and in the retort itself. The coal is carbonized in batches and the resulting coke is discharged after each period of carbonization. It was common practice in the early 1900s for manufactured gas plants to bury some of their wastes in unlined pits on site. Most of the coal tars, approximately 76 percent, were sold to coal tar refineries for further processing.

3.3 PAST INVESTIGATIONS

On April 8, 1987, Daniel Vilson, Environmental Sanitation Specialist for the St. Louis Division of Health, collected oil samples in the basement of the former Mound Street electric power plant, where hydraulic oil from electrical transformer allegedly was stored. Six samples were collected and analyzed for polychlorinated biphenyls (PCBs). According to this analytical report, none of the samples showed PCB contamination, though the detection limits were not included on the data transmittal (E & E/FIT 1988).

On September 17, 1987, the E & E/FIT conducted a site reconnaissance to aid in preparing a Preliminary Assessment report of the former Mound Street Power Plant. Six liquid samples were collected from the basement of the facility, and two samples were collected from two different manholes adjacent to the facility. All samples were screened for PCBs at a 1 part per million (ppm) detection limit. No PCB contaminants were identified by the Tracor gas chromatograph utilized by the FIT Field Analytical Support Program (FASP) (E & E/FIT 1988).

The E & E/FIT conducted a second site reconnaissance on November 20, 1990, to facilitat preparation of the work plan. The FIT observed seepage from the foundation and piping system of an abandoned pump house that was formerly owned by the Mound Street Power Plant. The pipes, which originate from the plant, had been plugged with concrete, but seepage was still leaching through the concrete. Because this pump house is located on the east side of the flood control levee, this leachate was observed to be seeping directly into the Mississippi River (E & E/FIT 1990).

SECTION 4. VASTE CHARACTERISTICS

4.1 VASTE STREAMS

The two waste products of primary concern are tar sludges (coal tars) and spent oxides. Ammonia wastes are by-products of this production process, but are not considered hazardous. Coal tar wastes are primarily polynuclear aromatic hydrocarbons (PAHs) and phenolics produced during coal or coke combustion and during the oil injection process. Some of the PAH compounds likely to be present in the tar wastes are carcinogenic and listed as RCRA Part 261 hazardous wastes. All PAHs can be considered as carcinogenic as benzo(a)pyrene, a Class A carcinogen. Spent iron oxide wastes are produced during the gas purification process where impurities are removed from the manufactured gas. Iron oxide wastes contain sulfur compounds, cyanide compounds, and small quantities of coal tar. Light aromatics such as benzene, toluene, and xylene (volatile organic compounds) also are occasionally constituents of coal tar wastes.

Two types of cyanide may be present at a coal gasification site: simple and complex cyanides. Simple cyanides are formed when cyanide reacts with an alkali or metal, producing a soluble material that can liberate a CN anion in water. Simple cyanides can be decomposed by bacteria in the soil (Environmental Research 1984). Complex cyanides are alkali-metal cyanides that are relatively insoluble (Environmental Research 1984). Complex cyanides, particularly the ferrocyanide compounds, are more resistant to biodegradation. These materials are associated with oxide wastes.

The trace metals most likely to be found on a coal gasification site are: arsenic, chromium, copper, iron, lead, nickel, and zinc (Environmental Research 1984). All are readily adsorbed onto soil particles. The mobility of these constituents is controlled by the pH of the soil. As a general rule, the solubility of these metals increases as pH decreases. Low pH values also reduce the cation exchange capacity of the soil matrix due to the preferential adsorption of H⁺ ions. Cation exchange is generally considered the major barrier to metal transport in soils. The strong tendency of metals to be bound to soil particles and organic matter limits their impact on ground water resources.

PAH and phenolic compounds may enter the atmosphere through

volatilization. Once in this matrix, the materials may undergo molecular or advective diffusion. (All further references to dispersion characteristics will infer both molecular and advective processes). PAH compounds are likely to undergo dispersion when introduced into surface water. If this occurs, the contaminants are very susceptible to adsorption onto clay particles suspended in the water. Depending on the nature of the surface water, this material may also volatilize; thus entering the atmosphere. Once in the surface water, the PAH compounds are prone to chemical alteration through biodegradation or photolysis. They are not readily absorbed to clay particles. These compounds may also undergo volatilization and limited biodegradation in surface water (Environmental Research 1984).

PAHs in ground water are also likely to undergo dispersion and adsorption processes. However, biodegradation of these materials is unlikely in this matrix (Environmental Research 1984). Phenolic compounds in ground water can be transported through dispersion. It is possible that these chemicals may undergo limited biodegradation in ground water environments.

In the soil matrix, PAHs can be involved in the adsorption process as well as biodegradation reactions. These materials may also undergo volatilization, leaching, and photolysis depending on site-specific characteristics. Phenolic compounds in the soil environment can be leached readily or removed through biodegradation (Environmental Research 1984).

PAH compounds are stable and tend to be retained in sediments. The specific stability of a particular PAH compound is dependent on its chemical structure (Environmental Research 1984; EPA 1980). Generally, the stability/solubility is inversely related to the molecular weight of the PAH. The arrangement of rings is also important. For example, anthracene is relatively soluble. It is a medium mass PAH composed of three linear rings. The arrangement of the rings allow this relatively massive molecule to be soluble. Benzo(a)pyrene is composed of a single ring surrounded by rings of three sides of its six sides. It is one of the more massive PAHs. The basic structures of the major PAHs are shown on Figure 4-1.

	Component	Formula	Structure	Boiling Point, C
	Benzene	C6H6	© ,	80
	Toluene	C-Hg	, ©	111
	Xylenes	C8H10	⊕ •• -	138-144
	Phenoi	C6H5OH	 6	181
	Cresois	C7H7OH	6 - . -	191-202
	Xylenois	C8H9OII	Q	201 - 227
_	Pyridine	CSHSN	Ø	115
	Naphtha i ene	C10H8	90	218
	Methylnaphthalenes	C11H10	•	241-245
	Dimethylnaphthalenes	C ₁₂ H ₁₂	- .	262-269
	Acenaphthene	C ₁₂ H ₁₀	•	277
	Carbazole	C ₁₂ H ₉ N	0,0	3 5 \$
	Fluorene	C13H10	600	297
	Anthracene	C14 ¹¹ 10	999	340
N S	Phenanthrene	C14H10	ම්ව න	340
	Fluoranthene	C16H10		393
NO.	Pyrene	C16 ¹¹ 10		394
PAH COMPOUNDS	Chrysene	C181112	0 00	436
۾ ا	Benz (a) anthracene	C ₁₈ ii ₁₂	000° a a	438
	Benzo(j)fluoranthene	C20H12	600	~480
[] []	Benzo(k)fluoranthene	C20 ¹¹ 12	96 <u>00</u>	480
	Benzo (a) pyrene	C ₂₀ H ₁₂	60 000	496
	Benzo (e) pyrene	C ₂₀ H _{1 2}		493
	Perylene	C ₂₀ H _{1 2}		460
Π^{\dagger}	Benzo(g,h,i)perylene	C ₂₂ !1 ₁₂	(0) (1) (1) (1) (1) (1) (1) (1) (1) (1) (1	500
	Benzo(b)chrysene	C22 ¹¹ 14	000	~500
	Oibenz(a,h)anthracene		0000 0000	•

Source: Environmental Research 1984

Chemical Compounds Associated with Coal Gasification

FIGURE 4-1

4.2 VASTE QUANTITY

The Laclede Coal Gas facility operated for almost 60 years, producing approximately 932,640 million cubic feet of gas (Andrew 1890-1940). According to the Handbook on Manufactured Gas Plants, approximately 1 gallon of tar waste is accumulated for every 1000 cubic feet of gas produced (Environmental Research 1984). Using this conversion factor, approximately 932 million gallons of tar waste were produced at the Laclede Coal Gas facility. According to the Handbook on Manufactured Gas Plants, approximatley 76% of the tar waste produced was sold. The other 24% was probably buried on site. Therefore, approximately 223,680,000 gallons of coal tar could be buried on site. It was common practice for FMGPs to bury their waste on site in pits or unlined trenches.

SECTION 5: PHYSICAL AND CULTURAL SETTING

5.1 CLIMATOLOGY AND DEMOGRPHY

The climate of the St. Louis, Missouri, area is classified as consistent with cold winters and long, hot summers. Heavy rains occur mainly in the spring and early summer months (USDA 1982). The total annual precipitation for St. Louis is 36 inches (USDA 1982). Net precipitation is 0 inches and the one-year, 24-hour rainfall is approximately 3.0 inches (USDA 1982). The average relative humidity in mid-afternoon is about 60 percent; with a higher humidity at night and the average at dawn is about 80 percent. The prevailing wind is from the south (USDA 1982).

The total population within four miles of the site is 207,102 (EPA 1989). The nearest individual is located approximately 1/4 mile from the site. The only resource within one-half mile of the site is the Mississippi River, located directly east of the site.

According to the Missouri Department of Conservation, it is not likely that any sensitive environments or critical habitats occur within a four-mile radius of the site. The possible exception is the Sicklefin Chub, (<u>Bybopsis meek</u>), which occurred in the Mississippi River in 1944 and may still occur in the area (Dickneite 1991).

5.2 TOPOGRAPHY AND DRAINAGE

The site topography is essentially flat with a very gentle slope (0 to 3 percent) to the east. Locally, the slope has been modified around buildings and other facilities. All runoff is impeded by a 500-year Corps of Engineers flood wall, which separates the site from the Mississippi River (Borgestede 1990). Overland flow is diverted into a storm sewer system that is connected to a sanitary sewer line. All water entering the sanitary sewer system flows underground to the Bissle Point Treatment Plant located approximately 2.5 miles upstream from the site. The water is treated by primary settling tanks; however, construction of a secondary treatment plant is three years underway. Treated water from this plant is discharged into the Mississippi River approximately 2.5 miles north (upstream) of the site (Borgestede 1990).

5.3 SOILS

The site is situated on Urban bottomland, which is defined as land with more than 85 percent of the surface covered by asphalt, buildings, or other impervious materials (USDA 1982). The area was built up to protect the site from flooding. The amount of fill in the area can range from 0 to over 20 feet. Because of this high degree of variability, classifying the exact nature of the area soils is impractical without a detailed on-site soil analysis. However, generally the soil types found at the Laclede Coal Gas site are periodologically characterized as loessial. These unconsolidated loess deposits of Quaternary Age can vary in thickness up to 50 feet. Residual soils underlie these loess deposits and were formed from decomposition of limestone bedrock (MGS 1971).

5.4 STRATIGRAPHY AND GROUND VATER

The upper bedrock beneath St. Louis County in east-central Missouri belongs to the Upper Mississippian and Lower Pennsylvanian systems. Figure 5-1 shows that these systems are subdivided, in descending order, into the Pleasanton, Marmaton, and Cherokee groups of the Pennsylvanian System, and the Mermacian Series of the Mississippian System (MGS 1961). Figure 5-2 is a generalized stratigraphic column primarily prepared from two well logs obtained from the Division of Geology and Land Survey, Missouri Department of Natural Resources (MDNR) and the drilling and sampling logs prepared during the SSI. The drilling and sampling logs are included in Appendix G.

According to the two well logs obtained from the MDNR and the drilling and sampling logs prepared during the SSI, the depth to bedrock varies in the site vicinity. The drilling and sampling logs (Appendix G) indicate that this depth ranges from 23 feet (at borehole 24) to 29 feet (at borehole 07). The well logs (Appendix F) indicate that the depth to bedrock is between 20 and 25 feet. However, this depth would increase as the distance to the Mississippi River decreases. To simplify the generalized stratigraphic column, the depth to bedrock was rounded to 30 feet. This 30 foot thickness is comprised of unconsolidated loess and alluvial (Mississippi River) deposits of the Quaternary System. Figure 5-3 depicts the alluvium thickness along the Mississippi River. In addition to these deposits, fill materials ranging from 0 to 20 feet or

PLEISTOCENE ALLEGOND SERIES ALLEGOND UPPRANTO /U PRANTO /U PRANTO /U	LEGEND
DESMOINE SIAN SERIES CYCNIAN SERIES CYCNIAN SERIES CYCNIAN CANONE CANO	MERAMECIAN SERIES [FAST-CENTRAL MISSOURI]
LACLEDE COAL	GAS SITE
ST. LOUIS, M	issouri

WASTE SITE TRACKING NO.: MO0579 PREPARED BY: JOHN C. PARKS

ECOLOGY AND ENVIRONMENT FIT DEC. 1987 SOURCE: MGSWR 1981

FIGURE 5-1: GENERALIZED STRATIGRAPHIC SECTION ST.LOUIS COUNTY, MISSOURL

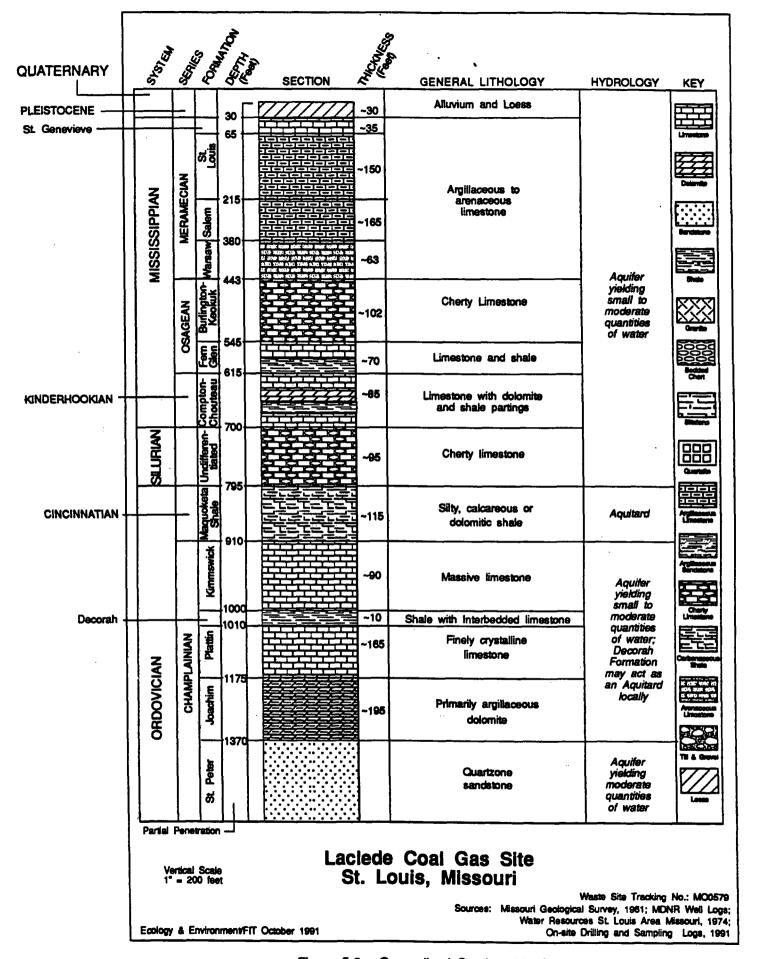
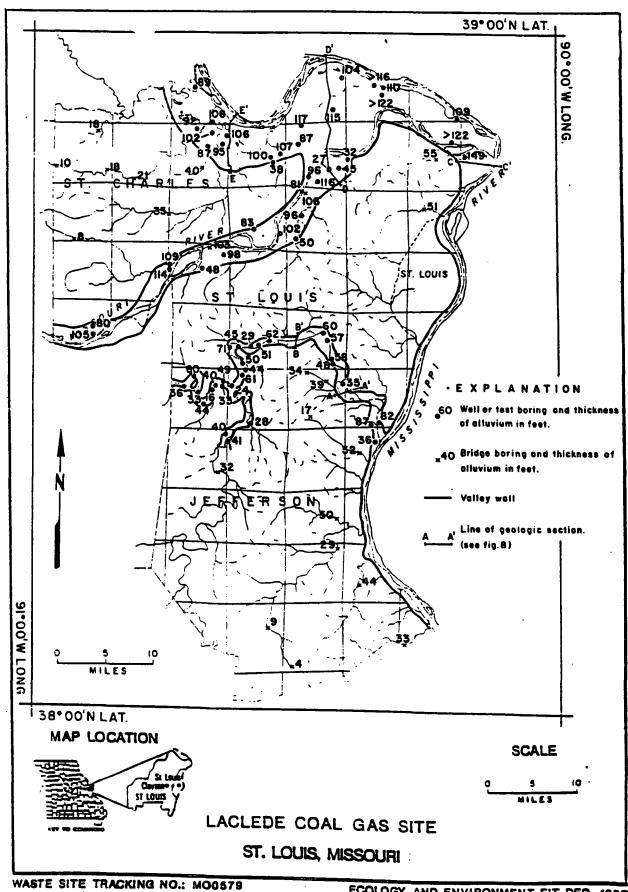


Figure 5-2: Generalized Stratigraphic Column



WASTE SITE TRACKING NO.: MO0579
PREPARED BY: JOHN C. PARKS

ECOLOGY AND ENVIRONMENT FIT DEC. 1987

Source: MGSWR 1974

FIGURE 5-3: ALLUVIUM THICKNESS ALONG THE MISSOURI. MISSISSIPPI AND MERAMEC RIVERS ST. LOUIS COUNTY MISSOURI

more was placed at the site in order to protect the area from flooding.

The first bedrock encountered below the site is the Mermacian Series of the Mississippian System. The Mermacian Series consists of four formations (Fm.), which in descending order include the St. Genevieve, the St. Louis, the Salem, and the Warsaw. These formations are composed primarily of limestone and have a total thickness of 300 to 450 feet in east-central Missouri (MDNR Undated).

The St. Genevieve Fm. consists of white, massively bedded, sandy, clastic limestone. Although there are some thin beds of finely crystalline limestone, the limestone is primarily coarsely crystalline and colitic. The lower portion of the formation is sandy, white to light tan or light clive gray, and is cross bedded and ripple marked. The middle of the formation consists of layers of red and gray chert with lenses and beds of sandstone occurring locally. Certain beds contain notable amounts of limonite which lines small cavities in the rock. Fossils are irregularly distributed in the formation. The thickness of this formation is approximately 30 feet in the county and approximately 35 feet at the site (MGS 1961; MDNR Undated).

The St. Louis Fm. consists of gray lithographic to finely crystalline, medium to massively bedded limestone which can be more than 100 feet thick in the county (MGS 1961). At the site, this formation is approximately 150 feet thick (MDNR Undated). Limestone brecia with a shale matrix between blocks of breccia is common in the lower portion of the formation. Blue to bluish-gray shale also occurs in thin beds throughout the formation. Chert is not common, but when observed it is usually brown and in the form of small angular fragments. Parts of the formation are locally dolomitic (MGS 1961).

The Salem Fm. thickest exposures are found in St. Louis, St. Genevieve, and eastern Perry Counties. The thickness of this formation generally ranges from 100 to 160 feet (MGS 1961). The thickness of this formation at the site is approximately 165 feet (MDNR Undated). The limestone contained in the formation is more dolomitic as compared to the other three formations in the Mermacian Series. The upper portion is fossiliferous and contains blastoid, crinoid, echinoid, and bryozoan debris. Also the upper portion contains speckled gray and tan chert (MGS 1961).

The Warsaw Fm. can outcrop in S* Louis County and where outcrops occur, the thickness of the formation is between 80 to 100 feet (MGS 1961). At the site, this formation is approximately 63 feet thick (MDNR Undated). This formation consists primarily of finely crystalline shaley, very fossilferous, dolomitic limestone in the lower portion, and dark, fissile shale in the upper portion (MGS 1961).

According to the MDNR well logs, the lower bedrock units beneath the site vicinity include: the Burlington-Keokuk Limestone Fm., the Fern Glen Fm., and the Compton-Chouteau Fm. of the Mississippi System; an undifferentialed fm. of the Silurian System; and the Maquoketa Shale Fm., the Kimmswick Fm., the Decorah Fm., the Plattin Fm., the Joachim Dolomite Fm., and the St. Peter Sandstone Fm. of the Ordovician System (MDNR Undated). These bedrock formations will be discussed briefly below.

The Burlington-Keokuk Fm., consists primarily of white to light buff, very coarsely crystalline, fossiliferous, crinoidal limestone. Layers of chert nodules are common, particularly in the upper portion. In east-central Missouri, the thickness of the combined formation is about 125 feet (MGS 1961). Near the site the combined formation is approximately 102 feet thick (MDNR Undated).

The Fern Glen Fm. consist of gray, grayish-green, and red limestone, and green and red calcareous shale. In central St. Louis County, the predominate color of the limestone is red. The lower portion is noncherty, whereas the upper portion contains layers of grayish chert. The total thickness of the formation in the county ranges from 20 to 45 feet (MGS 1961). In the site vicinity, the formation is approximately 70 feet thick (MDNR Undated).

The Compton-Chouteau Fm. consists of finely crystalline to sublithographic, finely bedded crinoidal limestone. The beds are separated by green shale partings. Some portions of the formation are dolomitic and contains bluish-gray to bluish-black chert. In east-central Missouri, this formation merges with the overlying Sedalia Formation and the combined formation is referred to as the undifferentialated Chouteau. The formation is relatively thin in the thickness in east-central Missouri, ranging from 6 to 12 feet (MGS 1961). However, near the site vicinity, the formation is approximately 85 feet thick (MDNR Undated).

Beneath the Mississippian-age formations: the site vicinity is a layer of undifferential Silurian-aged rocks. This unit is approximately 95 feet thick in the site area and consists primarily of cherty limestone (MDNR Undated).

Below the undifferential Silurian-aged rocks are the formations of the Ordovician System. The first formation, encountered below the site in the Ordovician System is the Maquoketa Shale Formation (MDNR Undated). This formation consists of thinly laminated, silty, calcareous or dolomitic shale with nodular and shaly lenses of limestone. The thickness ranges from 30 to 140 feet (MGS 1961). Near the site vicinity, the formation is approximately 115 feet (MDNR Undated). The next formation encountered below the site is the Kimmswick Formation. This formation consists of coarsely crystalline, white to light gray, medium bedded to massive limestone. Chert is irrregularly scattered throughout the formation. The foramtion ranges from 50 to 150 feet thick (MGS 1961). In the site vicinity, the formation is approximately 90 feet thick (MDNR Undated).

The Decorah Fm. lies below the Kimmswick formation in the site vicinity and consists of green and brown shales and interbedded limestone. The upper portion contains medium to thinly bedded, fossiliferous limestone with thin shale partings. The lower portion contains more shale with numerous, thin interbedded limestone. The formation varies in thickness from a few feet to more than 40 feet (MGS 1961). Near the site, the formation thickness is approximately 10 feet (MDNR Undated).

The Plattin Fm. underlies the Decorah Fm. near the site and consists of evenly bedded, dark gray, finely crystalline to sublithographic limestone. The formation also contains minor amounts of intercalated shale, and white chert nodules and layers are present throughout most of the formation. All of the formation in east-central Missouri is composed of dolomite. The formation can be as much as 450 feet thick north and west of Cape Girardeau County (MGS 1961). Near the site, the approximate thickness is 175 feet (MDNR Undated).

The Joachim Fm. lies below the Plattin Fm. in the site area and consists of yellowish-brown, argillaceaous dolomite which contains interbedded limestone and shale in this lower part (MGS 1961). Near the

site, the thickness of this formation is approximately 1/5 feet (MDNR Undated).

Below the Joachim Fm. in the site area, and the last formation to be discussed in this report is the St. Peter Sandstone Formation. This formation consist primarily of quartzose sandstone. The thickness of the formation is variable, ranging from less than 10 feet to more than 100 feet (MGS 1961).

Because there is an abundance of potable surface water in the site vicinity, ground water is not utilized as a source of drinking water. Generally, the bedrock aquifers of the region yield very small quantities of water; roughly 0 to 50 gallons per minute (gpm). The alluvial aquifers of Quaternary Age that lie along the Meramac, Mississippi, and the Missouri Rivers have much greater yields. Ground water flow in the area is generally to the south and east. The depth to the water table varies because of the variability in the depth of fill material. Generally, the water table is about two feet above the Mississippi River (Vandike 1990).

The bedrock aquifers in the region are divided into five discrete units, or groups. Group one, the Post-Maquoketa group, includes the strata above the Kimmswick Formation to the surface. Below this aquifer group lies the Maquoketa Shale, which is believed to act as an aquitard. Group two is the Ordovician-age Kimmswick-Joachim aquifer. Near the top of this unit is the Decorah Formation, which probably acts as a confining bed composed of shales and interbedded limestones. The remaining lower three aquifers are, in descending order: the St. Peter-Everton, Powell-Gascondade, and the Eminence-Lamotte. These units are separated primarily on the basis of unconformities, and are likely to be hydraulically connected. These deeper bedrock aquifers can be seen in Figure 5-4 and 5-5, which depict generalized hydrogeology of St. Louis County.

The migration of coal tar in ground water has been observed in several former coal gas manufacturing sites (Environmental Research 1984, E & E/FIT 1987). Coal tar is more dense than water and tends to migrate downward through porous material to a confining layer of less porous material. In areas where this behavior is exhibited, the following stratification (from top to bottom) may be expected: ground water with

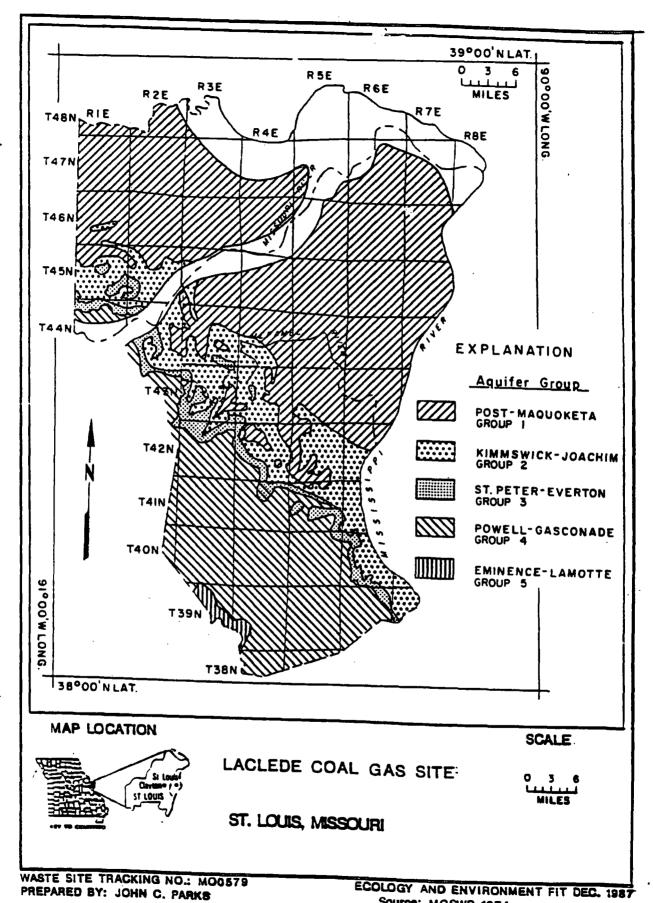
System	Series	Group	Formation	Aquifes Zroup	Thick- nass (fest)	Dominant lithology	Water-bearing character
	Noiocene	۱ اــــــــــــــــــــــــــــــــــــ	Aliuvium <u>i</u> /		v-150	Sand, gravel, silt,	Some wells vield more than 2,000 gpm.
unternary	Pleistocene		Loss Glacial till		0-110 0-55	Silt	Essentially not water yielding
1	Missourian	Pleasanton	Undifferentiated		0-75 0-90	Sheles, siltstones, "dirty" sandstones,	Generally yields very small quantities of
ennsylvantan	Desmoinesian	Cherokee	Undifferentiated	1 1	0-200	coal beds and thin	vater to welle.
,	Atokan		Undifferentiated			limestone beds.	Yields range from 0-10 gpm.
		†	Ste. Geneviewe Formation		0-160	Argillaceous to	
	Heramocian	·	St. Louis Limestons	1 1	0-150	Aretaceous limestess.	, 1
	İ	ļ	Sales Forestion	1 7	0-180	_	
		 	Vargaw Pormation Burlington-Keokuk Limestone	1	0-110 0-240	Charty limestone	-
Mississippiem	Congens		Fern Glea Formation	'	0-105	Red limestone and shale	lyields small to moderate quantities of water to
	Kinderhookias	Chouteau	· Undifferentiated		0-122	Limestone, dolomitic limestone, shale, and siltstone,	veils. Yields range from 5 to 50 gpm. Higher yields are
Devonisa	Upper	Sulpher Springs	Bushberg Sendstone Glen Park Linestone	- 1	0-60	Limestone and sandstone	
	<u> </u>		Grassy Creek Shale	1 '	0-50	Fissile, carbonsceous shale.	
Silurian	<u> </u>		Undifferentiated] '	0-200	Cherty limestone.	
			Maquoketa Shale		0-163	Silty, calcareous or dolomitic shale.	Probably constitutes a confining influence of water powerent.
	Cincinnatian		Cape Limestone Kimmerick	ــــــــــــــــــــــــــــــــــــــ	0-5 0-145	Argillaceous limestone.	니
		}	Formation	!	V-143	Maselve libestope	1
			Decorat Forestion	1	0-50	Shale with interbedded limestone.	quencities of water (
	Champlainian		Plattim Formation] 2	0-240	Finely crystalline limestone.	velle. Yields range from 3 to 50 gpm.
			Rock Laves Forestion	력 -	0-93	Dolomite and limestone,	probably acts as a
Ordovician			Joachim Dolemite St. Pater Sandstone	 	0-135	Primarily Argillaceous dolomice.	confining bed lecally
DE ROATE 1788	i		St. Pater Samustine	4	0-100	- Silty sendstone, cherr	y Yields moderate quanti-
			Everton Pormetion	3	0-130	limestone grading upward into quartzone sandstone.	ties of water to well Yields range from
			Powell Dolomite		0-150		Yields small to
	Canadian	1	Cotter Dolosite Jefferson City	-	0-320	Sandy and charty dolomites and	large quantities of water to wells.
	1		Dolomite	4		sandstone.	Yields range from 10
			Roubidoux Formation	Д	3-177	_	to 300 gpm. Upper
			Gasconade Dolonita Gunter Sandston Hember	10	0-380]	part of aquifer group yields only small amounts of water to wells.
		1	Inimence Dolomite Potosi Dolomite		J-17: J-325	Chefty dolomites, salt	Yields moderate to
Cambrian	Upper	Elvins	Derby-Doerun Colomice],	0-165	stones, sangstone, and shale.	vater to wells.
			Pavis Formation Econeterre purmatio	- -	2-3-385	_	10 to 400 gpm.
	<u> </u>		Lanotte Sandstone	===	235+	-	·
Precemorian	İ			i		Igneous and metemorphis rocks.	to wells in this ster
	<u> </u>	_1	i				

LACLEDE COAL GAS SITE ST. LOUIS, MISSOURI

WASTE SITE TRACKING NO.: MOQ579
PREPARED BY: JOHN C. PARKS

ECOLOGY AND ENVIRONMENT FIT DEC. 1987

Source: MGSWR 1974



Source: MGSWR 1974

FIGURE 5-5 : MAJOR AQUIFER DISTRIBUTION ST. LOUIS COUNTY MISSOURI-

dissolved organics; ground water with trapped coal tar; and below the confining layer, ground water with dissolved organics.

5.5 TARGETS

The drinking water needs of the city and surrounding community are provided primarily through the withdrawal of surface water from intakes on the Missouri, Mississippi, and Meramac Rivers. Withdrawal from these rivers total nearly 1.12 by 109 gallons per day (City of St. Louis. Undated). The municipal water intakes for the city of St. Louis and surrounding communities are approximately nine miles upstream from the site (MGS 1974). The combined flow from the Missouri and Mississippi Rivers averages approximately 1.12 by 10¹⁰ gallons per day. The Meramac has an average flow of 1.92 x 109 gallons per day. The municipal intake for the community of Illinois Metro East is located at mile marker 180.8, almost directly across the Mississippi River, about 1/4 mile east of the site (Figure 2-1). About 300,000 residents are served by this water supply. Illinois American Water Company draws approximately 15 to 20 million gallons per day from this intake during the winter months; withdrawal rates in the summary are much higher (Schlosser 1990). This is the only drinking water intake within 15 miles downstream of the site. Various industries have intakes on the Mississippi River, but their use is industrial.

According to the Missouri Department of Conservation, it is not likely that any sensitive environment or critical habitats occur within 15 miles downstream of the site. The possible exception is the Sicklefin Chub, (Hybopsis meek), which occurred in the Mississippi River in 1944 and may still occur in the area (Dickneite 1991).

SECTION 6: FIELD ACTIVITIES

Field work was conducted March 3 through 9, 1991. The sample series assigned to this activity was DSX44. The FIT members and their respective tasks were: Keith Brown, Team Leader; Anne Melia, FASP coordinator; Chris Williams, Geoprobe operator and sampler; Patty Roberts, Assistant Geoprobe operator and sampler; Wes McCall, drilling supervisor and Site Safety Officer; John Peck, sampler and health and safety monitor; and Jon Strobel, sampler and assistant FASP coordinator. The FIT contracted John Mathes & Associates, Inc. to conduct all subsurface drilling. On-site personnel for John Mathes & Associates, Inc. were: Keith Bunselmyer, driller; and Jeff Crank and Jim Burker, driller's assistants. Additionally, William Oberle and Jacalyn Wheeler, E & E/FIT, were present on site March 6 and 7, 1991, to conduct a health and safety, and technical field audits.

The primary chemical hazards associated with the Laclede Coal Gas site involve soils contaminated with cyanide salts and/or PAHs. These materials could pose inhalation, direct contact, and ingestion hazards. Samples were collected in level-D and level-C personal protection. An HNu photo-ionization detector with 10.2 eV probe was used to monitor ambient levels of volatile compounds in the breathing zone. If the HNu registered readings above predetermined action levels, personal protection was upgraded to level-C. Otherwise, samples were collected in level-D personal protection.

6.1 SUBSURFACE SOIL SAMPLING

Subsurface soil sampling was conducted March 4 through 9, in an attempt to determine the areal and vertical extent of subsurface contamination. Initially, 24 locations were drilled with solid stem augers (Figure 6-1; Table 6-1). Their depths ranged from 2 feet to 38 feet. Originally, the work plan called for 18 locations to be drilled with solid stem augers to approximately 20 foot depths, with the exception of 1 location which was to be drilled to 50 feet deep. This deep sample was intended to determine the depth to bedrock and help assess the vertical contaminant zone. The 6 additional locations were added because subsurface rubble prohibited the advancement of the augers

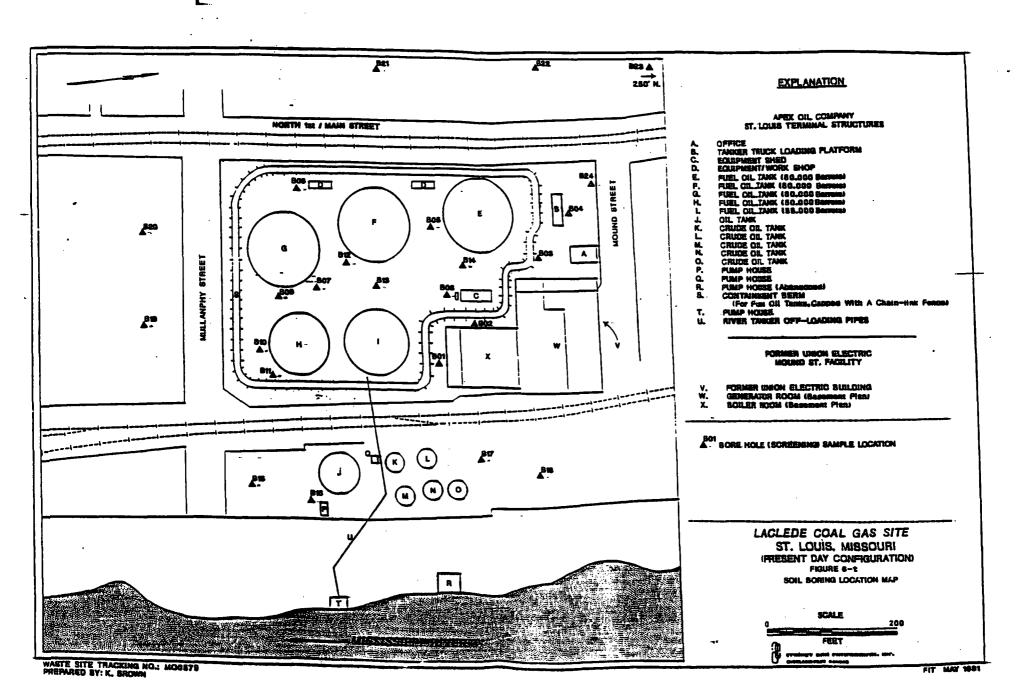


Table 6-1
Borehole Summary
Laclede Coal Gas
St. Louis, Missouri
Sample Series DSX44
March 1991

======================================	Depth (ft)	Approximate Location
BO1	18	South of former Mound Street Power Plant
BO2	21	Southwest of former Mound Street Power Plant
в03	19	South of P, F, & T Office
BO4	19	North of tanker truck loading platform
B05	26	Approximately 15 feet south of northern-most fuel
		oil tank
В06	38	North of center of site within containment berm
B07	29	South of center of site within containment berm
в08		Southwest corner of site within containment berm
В09	<u></u>	South of center of site within containment berm
B10		Southeast corner of site within containment berm
B11	15	Southeast corner of site within containment berm
B12	11	Center of site within containment berm
B13	10	Center of site within containment berm
B14	26	North center of site within containment berm
B15	31	Southeast corner of site
B16		Southeast corner of site approximately 100 feet
		north of B15
B17	33	Northeast corner of site
B18	33	Northeast corner of site approximately 150 north
		of B17
В19	30	South of Mullanphy Street
B20	30	South of Mullanphy Street approximately 200 feet
		north of B19
B21		Approximately 80 feet north of North 1st Street
B22 ·		Approximately 300 feet north of B21
B23		Approximately 450 feet north of B22
B24	23	Northeast corner of site

Note: See Figure 6-1 for borehole locations.

at several locations to only a few feet and, therefore, a representative sample could not be collected.

Samples were collected off the auger flights and screened on site for PAHs and VOCs in the FASP mobile laboratory (Tables 6-2 and 6-3). The samples were collected at 5 foot intervals: the PAH sample was a composite of 5 aliquots, 1 collected every foot; the VOC sample was a grab sample. Four of the sample locations determined by FASP analysis to be the most contaminated were resampled as follows: hollow stem augers equipped with continuous samplers were advanced, offset 1 to 5 feet from the screening boring (Figure 6-2, Table 6-4). A background location was sampled in the same manner. Due to subsurface rubble and debris, the continuous sampler could not be utilized at all locations. A split spoon sampler equipped with a 140 pound hammer was used at locations where the continuous sampler would not work. Samples were collected according to E & E, Inc. Standard Operation Procedures (SOPs) for borehole sampling, Gentech 5.9. Samples from the 4 locations that were resampled were submitted to EPA for volatiles, semi-volatiles, cyanides, and total metals analysis.

The background samples were inadvertently discarded along with the screening samples. This was not discovered by FIT until the SSI was completed and the team had returned to Kansas City. Split samples had been collected at all locations where samples were submitted to the EPA and which were also on PF & T property. Kathy Enright, of E & E/TAT in St. Louis, Missouri, sent the split samples collected for PF & T to E & E's Kansas City office. These samples were once again split, and 1 set was submitted to EPA to be used as the background sample. The other set was returned to PF & T.

The background samples were collected at a different location than stated in the work plan because bedrock was encountered at approximately 2 feet below the surface at this location; therefore, subsurface soil samples could not be collected there. FASP analysis had revealed that borehole #B24, the farthest upgradient screening sample, was below detection limits for all target compounds. It was known that the depth to bedrock was approximately 20 feet. Therefore, E & E/FIT relocated

C1- H	Benzene	Toluene	M-xylene	
Sample #	(µg/kg)	(µg/kg)	(µg/kg)	Depth (feet)
BO1 A	1,200	380	1,700	0 - 5
BO1 B	9,100	1,200	19,000	5 - 10
B01 C	18,000	710	65,000	10 - 15
B01 D	17,000	770	79,000	15 - 18
BO2 A	6,300	43,000	240,000	0 - 5
BO2 B	6,100	1,700	57,000	5 - 10
B02 C	69,000	110,000	570,000	10 - 15
BO2 D	7,500	650	33,000	15 - 21
BO3 A	1,040	22,000	22,000	0 - 5
BO3 B	310	11,000	53,000	5 - 10
BO3 C	1,800	6,300	3,500	10 - 15
BO3 D	860	9,000	5,300	15 - 19
BO4 A	< 250	450	5,600	0 - 5
BO4 B	310	480	5,900	5 - 10
B04 C	< 250	250	5,100	10 - 15
B04 D	< 250	ND	440	15 - 19
BO5 A	460	ND	ND .	0 - 5
BO5 B	ND	ND	ND	5 - 10
B05 C	ND	ND	420	10 - 15
B05 D	ND	ND	410	15 - 20
BO5 E	ND	ND	ND	20 - 26
B06 A	48,000	1,700	120,000	0 - 5
B06 B	20,000	1,800	68,000	8 - 10
BO6 BA	93,000	120,000	220,000	0 - 5
B06 BB	27,000	91,000	260,000	5 - 10
BO6 BC	ND	ND	ND	10 - 15
BO6 BD	< 250	< 250	ND	15 - 20
BO6 BE	12,000	14,000	54,000	20 - 25
BO6 BF	43,000	66,000	23,000	25 - 30
BO6 BG	20,000	16,000	26,000	30 - 35
B06 BH	29,000	43,000	94,000	35 - 38
B07 A	1,000,000	17,000	2,500,000	3 - 8
BO7 B	670,000	12,000	2,100,000	8 - 13
B07 C	370,000	ND	> 1,200,000	13 - 18
B07 D	140,000	2,500	> 21,000,000	18 - 23
BO7 E	22,000	490	120,000	23 - 26
B07 F	51,000	2,800	340,000	26 - 29
B10 B	< 250	410	ND	5 - 10
B11 A	ND	ND	ND	0 - 5

Table 6-2 (cont.)

	Benzene	Toluene	M-xylene	
Sample #	(µg/kg)	(µg/kg)	(µg/kg)	Depth (feet)
B11 B	ND	ND	ND	5 - 10
B11 C	ND	ND	ND	10 - 15
B12 A	< 250	< 250	1,500	0 - 5
B12 B	910	1,600	4,400	5 - 11
B13 A	570	420	< 250	0 - 5
B13 B	670	450	2,300	5 - 10
B14 A	270,000	11,000	> 160,000	2 - 8
B14 B	790	9,100	> 24,000	8 - 13
B14 C	400,000	14,000	> 160,000	13 - 18
B14 D	340,000	12,000	> 210,000	18 - 23
B14 E	310,000	8,900	> 195,000	23 - 26
B14 F	2,200,000	63,000	> 1,000,000	
B14 SS	93,000	110,000	440,000	
B15 A	680	490	> 29,000	0 - 5
B15 B	290	ND	10,000	5 - 10
B15 C	< 250	ND	7,400	10 - 15
B15 D	< 250	ND	2,700	15 - 20
B15 E	ND	ND	ND	20 - 25
B15 F	ND	ND	ND	25 - 31
B16 A	820	860	15,000	0 - 3
B17 A	540	ND	21,000	0 - 5
B17 B	ND	ND	2,700	5 - 10
B17 C	ND	ND	ND	10 - 15
B17 D	ND	ND	ND	15 - 20
B17 E	ND	ND	ND	20 - 25
B17 F	ND	ND	ND	25 - 30
B17 G	ND	ND	ND	30 - 33
B18 A	< 250	ND	ND	0 - 5
B18 B	ND	ND	ND	5 - 10
B18 C	ND	ND	ND	10 - 15
B18 D	ND	ND	ND	15 - 20
B18 E	ND	ND	ND	20 - 25
B18 F	ND	ND	ND	25 - 30
B18 G	ND	ND	ND	30 - 33
B19 A	ND	ND	ND	0 - 5
B19 B	ND	ND	ND	5 - 10
B19 C	ND	ND	ND	10 - 15
B19 D	ND	ND	ND	15 - 20
B19 E	ND	ND	ND	20 - 25
B19 F	ND	ND	ND	25 - 30
B20 A	320	ND	ND	0 - 5
B20 B	ND	ND	ND	5 - 10
B20 C	ND	ND	ND	10 - 15
B20 D	ND	ND	ND	15 - 20
				13 - 20
		1		
		<u> </u>		

Table 6-2 (cont.)

Sample #	Benzene (µg/kg)	Toluene (µg/kg)	M-xylene (μg/kg)	Depth (feet)
B20 E	ND	ND	ND	20 - 25
B20 F	ND	ND	ND	25 - 30
B22 A	ND	ND	ND	
B23 A	ND	ND	ND	0 - 4
B24 A	ND	ND ND	ND	0 - 5
B24 B	ND) ND	ND	5 - 10
B24 C	< 250	ND	ND	10 - 15
B24 D	400	ND	550	15 - 20
B24 E	460) ND	430	20 - 23
401	ND	, ND	ND	N/A
402	ND	ND	1,600	N/A
403	< 250	< 250	420	N/A
301	ND	ND	ND	N/A
302	ND	ND	ND	N/A
303	ND	ND	ND	N/A
304	ND	ND	ND	N/A
DW 1	188	77	1,000	N/A
DW 2	330	48	1,100	N/A
CDW	ND	ND	ND	N/A
SS-Rin-	ND	ND	ND	N/A
sate				1
		-		

Note: Detection limit for soil/sediment samples = 250 μ g/kg. Detection limit for water samples = 25 μ g/L. See Figures 6-1 and 6-2 for sample locations.

BO1 A through B24 E = soil samples

401 through 403 = sediment samples

301 through 304 = surface water samples

DW = decon water (units are µg/L)

CDW = clean decon water

i = no detection

N/A = not applicable

SS = split spoon

Table 6-3 FASP Screening Results for PAHs LaClede Coal Gas St. Louis, Missouri Sample Series DSX44 March 1991 Reported in µg/kg

Sample	Fluor-		Benzo(k)	Benzo(a)	
#	anthene	Pyrene	fluoranthene	pyrene	Comments
BO1 A	16,000	3,700	NA	NA	
B01 B	27,000	12,000	NA	NA	
B01 C	56,000	40,000	NA	NA	
BO1 D	13,000	5,200	NA	NA	
BO2 A	8,000	ND	ND	ND	
BO2 B	15,000	ND	ND	ND	
B02 C	ND	ND	ND	ND	
B02 D	ND	ND	ND	ND	
BO3 A	ND	ND	ND	ND	
BO3 B	ND	ND	ND	ND	1
B03 C	ND	ND	ND	ND	
B03 D	ND	ND ND	ND	ND	
B04 A	ND	ND ND	ND	ND	
B04 B	12,000	ND	ND	ND	
B04 C	9,000	ND	ND	ND	
B04 D	< 500	ND	ND	ND	chyrsene & benzo(a)
		•			anthracene = 1,200
BO5 A	ND	ND	ND	ND	
B05 B	ND	ND	ND	ND	
B05 C	ND	ND	ND	. ND	
B05 D	ND	ND	ND	ND	two unknown peaks
BO5 E	ND	ND	ND	ND	İ
B06 A	55,000	110,000	54,000	40,000	anaphthalene, fluorene, anthracene, chrysene
B06 B	Ble	v out GC. pea	ks too large t	' o integrate	
BO6 BA		ND	ND	I ND	I
BO6 BB	1	ND	ND	ND	
BO6 BC	I.	ND	32,000	11,000	other identified peaks
B06 BD	1	ND	ND ·	ND	Total Total Total Found
BO6 BE		ND	ND	ND	
BO6 BF		ND	ND	ND	
B06 BG		ND	ND	ND	Ì
B06 BH		ND	ND	ND	İ
B07 A	•	too large to	•	time to re	analvze
во7 в	160,000	50,000	50,000	70,000	other PAHs present
B07 C	130,000	130,000	33,000	76,000	other PAHs present
B07 D	12,000	9,600	3,700	3,900	P. C.
B07 E	3,300	3,800	3,500	3,000	other PAHs present
B07 F	20,000	19,000	4,200	8,000	other PAHs present
B10 B	14,000	16,000	7,800	9,100	other PAHs present
				1	F
			6-8		

6-8

Table 6-3 (cont.)

B11 A	2,400	1,600	4 200		T	
B11 B	?,100	2,100	4,200	ND 10 000	other	PAHs present
B11 C	7,800	•	7,000	10,000	other	PAHs present
B12 A	14,000	ND 12,000	6,000	12,000	Ì	
B12 B	38,000		14,000	15,000		
B13 A	150,000	29,000	12,000	15,000	Į	•
B13 B	12,000	73,000	46,000	44,000		
B13 C	32,000	10,000	10,000	12,000	'	
B14 A	1,600,000	34,000	28,000	32,000	ļ <u>.</u>	
B14 B	2,200,000	580,000	280,000	63,000		identified peak
B14 C	4,200,000	1,000,000	370,000	310,000		identified peak
B14 C	4,200,000	1,600,000	1,200,000	1		identified peak
B14 E	3,200,000	1,300,000	490,000	110 000		identified peak
B14 E		4,100,000	1,500,000	140,000	other	identified peak
B14 F	1 400 000	w out GC, pea	aks too large			
B14 33	, , ,	880,000	230,000	130,000	ļ	
	ND 4 000	ND 1 300	ND	ND		
B15 B	4,900	1,300	ND	ND	ļ	
B15 C	ND ·	ND	ND	ND	ļ	
B15 D	ND	ND	ND	ND	ļ	
B15 E	ND	ND	ND	ND		
B15 F	ND	ND	ND	ND	!	
B16 A	19,000	30,000	ND	ND	early	peaks
B17 A	ND	ND ND	ND	ND	1	
B17 B	ND	ND	ND	ND		
B17 C	ND	< 500	ND	ND]	
B17 D	ND	ND	ND	ND		
B17 E	ND	ND	ND	ND	ĺ	
B17 F	ND	ND	ND	ND		
B18 A	ND	ND	ND	ND		
B18 B	ND	ND	ND	ND	ļ	
B18 C	ND	ND	ND	ND		
B18 D	ND	ND	ND	ND	ļ	
B18 E	ND	ND	ND	ND		
B18 F	ND	ND	ND	ND		
B18 G	ND	ND	ND	ND	1	
B19 A	ND	ND	ND	ND		
B19 B	ND	CA C	ND .	ND	1	
B19 C	ND	ND	ND	ND		
B19 D	ND	ND	ND	ND		
B19 E	18,000	16,000	14,000	12,000		
B19 F	4,500	ND	2,800	ND	1	
B20 A	ND	2,900	< 500	ND		
B20 B	4,100	4,000	ND	545		
B20 C	ND	ND	ND	ND		
B20 D	ND	ND ND	ND	ND	1	
B20 E	ND	ND	ND	ND	1	
B20 F	2,600	3,200	3,500	4,500	1	
B22 A	3,800	3,500	2,800	ND		
	1	[1	
	L	<u> </u>			<u></u>	
			6-9			

Tabl ? 6-3 (cont.)

B23 A	ND	ND	ND	ND	
501					peaks close to reten-
İ					tion time, but do not
ļ	Ì	}			match
402	ND	ND	ND	ND	
403	2,600	5,400	4,200	3,800	•
DDW	ND	ND	ND	ND	no peaks (µg/L)
SS Rin	ND	ND	ND	ND	no peaks (µg/L)
sate	İ	İ			
B24 A	ND	ND	ND	ND	1
B24 B	ND	ND	ND	ND	1
B24 C	ND	ND	ND	ND	}
B24 D	ND	ND	ND	ND	1
B24 E	ИD	ИD	ND	ND	\
j	Ì	j			

Note: Sample locations are shown on Figures 6-1 and 6-2. Sample 501 is an oil sample collected from tank #7, shown on Figure 6-2. Detection limits for soil/sediment = $500 \mu g/kg$. Detection limits for water samples = $15 \mu g/l$.

DDW = dirty decon water

SS = split spoon

NA = not analyzed

ND = not detected

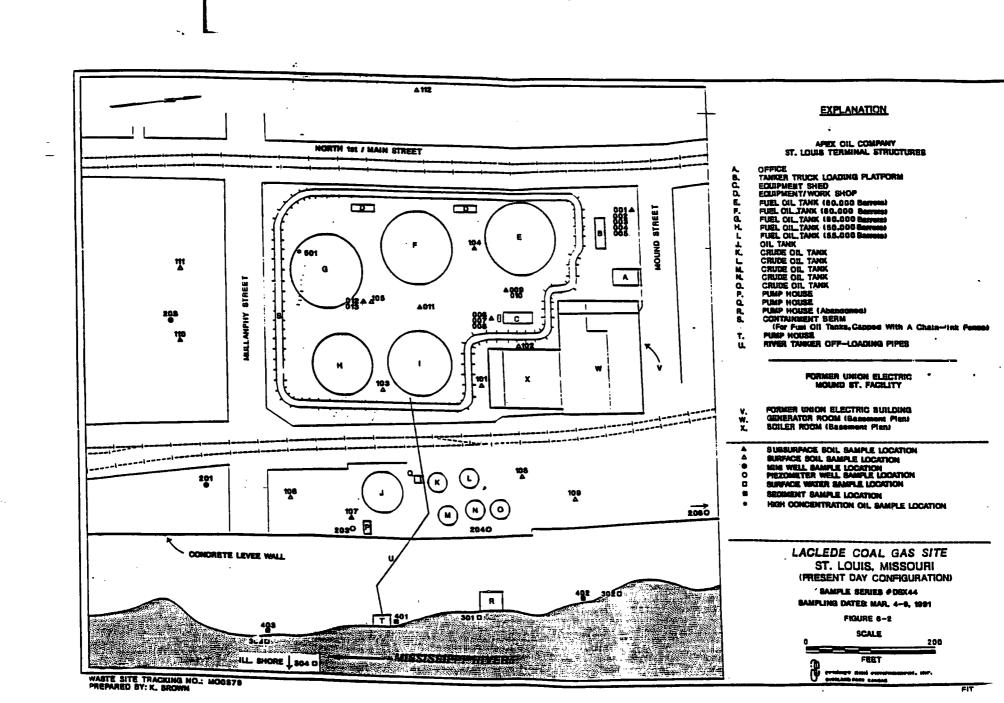


Table 6-4 CLP Soil/Sediment Sample Summary LaClede Coal Gas St. Louis, Missouri Sample Series DSX44 March 1991

Sample #	Depth (feet)	Location Description
001	3 - 7	5 feet northwest of B24
002	7 - 11	5 feet northwest of B24
003	11 - 15	5 feet northwest of B24
004	15 - 19	5 feet northwest of B24
005	19 - 23	5 feet northwest of B24
006	3 - 8	2.5 feet north of BO6
007	8 - 13	2.5 feet north of BO6
800	8 - 12	2 feet south of BO6
009	4 - 8	3 feet northwest of B14
010	12 - 14	3 feet northwest of B14
011	11 - 17	1.5 feet west of BO7
012	5 - 11	1 foot north of B13
013	11 - 12.5	1 foot north of B13
101	0 – 2	B01
102	0 – 2	B02
103	0 - 2	South central tank farm
104	0 - 2	B05
105	0 - 2	B07
106	0 - 2	B15
107	0 - 2	B16
108	0 - 2	B17
109	0 - 2	B18
110	0 - 2	B19
111	0 - 2	B20
112	0 - 2	B21
401	N/A	150 feet south of southeast corner
	İ	of the abandoned pump house
401D	N/A.	150 feet south of southeast corner
	ĺ	of the abandoned pump house
402	N/A	110 feet north . northeast corner
		of the abandoned pump house
403	N/A	330 feet south of southeast corner
		of the abandoned pump house

Note: Samples 001 through 013 collected from five aliquots, except VOCs were grab samples. Samples 101 through 112 collected from two aliquots. All samples were submitted to EPA for cyanide, semi-volatiles, and total metals analyses. Samples 001 through 113 and 401 through 403 were also analyzed for volatiles. Samples 401 through 403 were also analyzed for total petroleum hydrocarbons. See Figure 6-2 for sample locations.

D = Duplicate

N/A = Not applicable

the background sample location to the farthest northwest corner of PF & T property. This location is approximately 60 feet northwest (upg-adient) of borehole #B04 (Figure 6-1).

6.2 SURFACE SOIL SAMPLING

Surface soil sampling was conducted March 4 through 7, to determine whether contamination was present in the 0- to 2-foot interval. Twelve samples were collected (Figure 6-2, Table 6-4). According to the work plan, these samples were to be collected using a power auger. How ver, due to excessive amounts of gravel and rubble present, the power auger could not be used efficiently. Therefore, all but one surface soil sample was collected with the subcontractor's drill rig.

Each sample was collected off the auger flights and consisted of two aliquots collected from one- and two-foot depths. The sample not collected with the drill rig was obtained with the use of the power auger. Three of the 12 surface soil samples collected, 110, 111, and 112, were inadvertently discarded along with the screening samples. Keith Brown and Scott Hayes returned to the site on March 13 and 14 and recollected these samples with a power auger. All samples were submitted to EPA for semi-volatile, cyanide, and total metals analyses. Samples were collected in accordance with the Region VII E & E/FIT SOP for soil sample collection, Geotech 5.17.

6.3 GROUND WATER SAMPLING

Ground water sampling was conducted March 5 through 9, to determine whether any coal tar wastes present have leached into the shallow ground water or migrated off site via the ground water. A total of 9 samples, including a duplicate, field blank, rinsate, and extra volume sample were collected (Figure 6-2, Table 6-5). The work plan called for the installation of 4 temporary mini-wells with the Geoprobe. Due to subsurface rubble, only 2 mini-wells were installed. However, the Corps of Engineers (COE) has installed piezometer wells along the west side of the Mississippi River levee wall. The COE wells located just east of the site were sampled with the use of the Geoprobe vacuum system. One other COE piezometer well was sampled; it is located approximately 1,500 feet north (upgradient) of the site along the levee wall. This location

Table 6-5
Water Sampling Summary
LaClede Coal Gas
St. Louis, Missouri
Sample Series DSX44
March 1991

		Static Water	
Sample #	Depth (feet)	Level (SWL) (feet)	Location/Description
201	26	24	64 feet south from center of Mullanphy Street and 50 feet east of easternmost railroad
202	33	24	147 feet west of easternmost railroad and 78 feet south of center of Mullanphy Street
203	47.25	27.25	147 feet north of center of Mullanphy Street and 40 feet west of levee wall
204	48	26	363 feet north of center of Mullanphy Street and 25 feet west of levee wall
205F	N/A	N/A	Rinsate of augers and split spoon sampler
206	52	29	2,118 feet north of center of Mullanphy Street and 20 feet west of levee wall
206D			Duplicate of 206
207F	N/A	N/A	Field Blank
208	N/A	N/A	Geoprobe pipe rinsate sample
209F	N/A	N/A	Trip blank
301	Surface	N/A	Southeast corner of abandoned pump house
301D	Duplicat	e of above	
302	Surface	N/A	170 feet north of northeast corner of abandoned pump house
303	Surface	N/A	330 feet south of southeast corner of abandoned pump house
304	Surface	N/A	Illinois American drinking water intake in East St. Louis
304D			Duplicate of 304
501	Surface	N/A	Oil sample collected from PF & T tank #80-1

Note: See legend on next page.

Legend for Table 6-5

Note: Sample locations are shown on Figure 6-2. All water samples were submitted to EPA for volatiles, semi-volatiles, and total and dissolved metals analyses. Samples 301 through 304D were also analyzed for Total Petroleum Hydrocarbons. Sample #501 is an oil sample which was submitted to EPA for Total Petroleum Hydrocarbons, volatiles, and semi-volatiles analyses. Cyanide samples were preserved with sodium hydroxide. Metals samples were preserved with nitric acid.

D = duplicate
F = field blank or trip blank
N/A = not applicable

was chosen for the ground water background sample. The original back-ground location could not be sampled because bedrock was only 2 feet deep. All samples were collected following the E & E Region VII SOP for ground water collection, Gentech 5.11. Total and dissolved metals samples were preserved with nitric acid; cyanide samples were preserved with sodium hydroxide; and all samples were stored on ice. Samples were delivered to EPA for volatiles, semi-volatiles, cyanide, total metals, and dissolved metals analyses. The field parameters, temperature, pH, and conductivity were measured and recorded in the field (Table 6-6).

6.4 SURFACE VATER/SEDIMENT SAMPLING

Surface water/sediment sampling was conducted March 6 through 7 to determine whether contaminants have migrated via the ground water/surface water interface. A total of 6 surface water samples were collected, including 2 duplicates (Figure 6-2, Table 6-5). All samples were collected according to the EPA Region VII SOP for surface water sample collection, 2334.7A. Samples were submitted to EPA for Total Petroleum Hydrocarbons, volatiles, semi-volatiles, cyanide, total metals, and dissolved metals analyses. Total and dissolved metals samples were preserved with nitric acid, cyanide samples were preserved with sodium hydroxide, and all samples were stored on ice. The field parameters temperature, pH, and conductivity, were measured and recorded in the field (Table 6-6).

A total of 4 sediment samples were collected, including a duplicate (Figure 6-2, Table 6-4). All samples were collected according to the EPA Region VII SOP for sediment sample collection, 2334.8A. Samples were submitted to EPA for Total Petroleum Hydrocarbons, volatiles, semi-volatiles, cyanide, and total metals analyses. Surface water and sediment samples were screened on site in the mobile lab for the presence of PAHs and VOCs (Tables 6-2 and 6-3).

6.5 PRODUCT SAMPLING

One oil sample was to be collected from the basement of the former Mound Street Power Plant, if the building was standing and oil was present in the basement. At the time of the SSI, the building was in the process of being razed. Due to the danger involved with entering

Table 6-6
Field Parameters for Water Samples
LaClede Coal Gas
St. Louis, Missouri
Sample Series DSX44
March 1991

Sample #	Temperature (C°)	рН	Conductivity (umhos)
301	7	8.15	440
301D	7	8.15	440
302	6	7.87	450
303	6	7.27	360
304	3	7.05	350
304D	3	7.05	350
201	11	7.58	1,500
202	7	7.81	2,000
203	9	6.32	1,500
204	9	6.21	1,400
206	11	6.14	1,600
206D	11	6.14	1,600
207F			
208	4	7.00	16
209F			

Note: See Figure 6-2 for sample locations.

this building, the mple was not collected. However, an oil sample (sample 501) was collected from an 80,000-gallon tank on site (Figure 6-2). This tank contained #6 oil, the same type of oil that was spilled in 1986 and entered the basement of the former power plant. This oil sample was submitted to EPA for total petroleum hydrocarbons, VOCs, and semi-volatiles analyses. It was also screened on site in the mobile lab. The chromatogram peaks were close to the retention times for the contaminants being screened, but did not match (Table 6-3).

SECTION 7: ANALYTICAL RESULTS

7.1 SURFACE SOIL SAMPLES (0-2 FEET)

The surface soil (0-2 feet) sampling data indicated that the majority of the 7-acre site is significantly contaminated with cyanide at greater than three times the detection limit of the background sample 112 (Figure 6-2; Table 7-1). Analysis of off site samples 110, 111, and 112 revealed no cyanide contamination over their detection limits of 6.2, 6.2, and 6.7, respectively (Figure 6-2).

Surface soil PAH contamination was less extensive than the surface soil cyanide contamination. Four out of the 12 samples collected contained a total PAH content greater than five times the concentration of the background sample. These were samples 104, 105, 107 and 109 (Figure 6-2; Table 7-1).

The cyanide contamination found on site can be attributed to the FMGP based on the fact that cyanide is an oxide waste produced during the gas purification process (see Section 4 for details). There should not be any attribution from current on site activities, because cyanide is not a constituent normally found in oil or asphalt.

Attribution of PAH contamination in the surface soil samples is difficult to assess. The oil sample data indicated significant concentrations of many PAH compounds (Appendix D, Sample 501). PF & T currently stores #6 oil and asphalt on site in very large quantities. In 1986, PF & T experienced a spill of said oil which covered most of the site. Thus, surface soil PAH contamination could be attributed to PF & T activities or to the FMGP, based on the fact that it was common practice for FMGPs to bury tar and purifier waste on site. PAHs are a major constituent of coal tar waste.

7.2 SUBSURFACE SOIL SAMPLES

FASP screening results indicated that the highest PAH contamination zone was located in the central portion of the tank farm currently on site (Figure 6-1; Tables 6-2 and 6-3). The approved work plan allowed E & E/FIT to collect CLP subsurface soil samples at four locations. A total of eight samples (006-013) were collected at depths ranging from 3

Table 7-1
PAHs and Cyanide In Surface Soils
Laclede Coal Gas Site
St. Louis, Missouri
E & E/FIT; March 1991
Sample Series DSX44
(mg/kg)

Compound	101	102	103	104	105	106	107	108	109	110	111	112*
Napthalene							60	1.1	0.51			0.41U
2-Methylnaphthalene							13	2.5				0.41U
2-Chloronaphthalene												0.41U
Acenaphthylene								2.4	0.46			0.41U
Acenaphthene								0.69				0.41U
Fluorene					12			3.1				0.41U
Phenanthrene				20	55	1.5			2.2	1.5	0.85	0.44
Anthracene									0.78			0.41U
Pluoranthene				14	38	2.0			2.8	2.2	0.77	1.3
Pyrene		21	13	50	83	2.2			6.7	2.3	0.81	1.4
Benzo(a)anthracene				21	28	1.3			4.5	1.5	0.48	0.79
Chrysene				29	29	1.2			4.3	1.6	0.55	0.85
Benzo(b)fluoranthene			 .		17	1.0			4.9	1.1	0.44	0.61
Senzo(k)fluroanthene					12	0.89			3.4	1.2	0.45	0.68
enzo(a)pyrene				16	22	1.1			4.2	1.3	0.44	0.70
ndeno(1,2,3-CD) Pyrene			· 			0.66			2.7	0.76	.	0.41U
ibenzo(a,h)anthracene												0.41U
ibenzo(g,h,i)perylene					12	0.71			2.6	0.76		0.41U
otal PAH Content		21	13	150	308	12.5	73	9.8	40.0	14.2	4.8	6.8
yanide	33		94	220	190		14	98	35			6.7U

^{* =} Background sample
-- = Undetected (U); Detection limit stated for bckground sample.

Note: See Figure 6-2 for sample locations and Appendix D for complete analytical results.

to 17 feet. An additional five samples (001-005), are collected at a background location (Figure 6-2; Table 6-4). CLP data revealed significant PAH contamination at least five times greater than background concentration at all four sampling locations (Table 7-2). Cyanide contamination was found to be significant to a depth of 11+ feet (Table 7-2).

All the soil samples were also analyzed for total and dissolved metals, as well as volatiles. However, results proved to be negligible when compared to background concentrations with the exception of significant concentrations of benzene, toluene, ethyl benzene, and total xylenes (BTEX) in samples 006 through 011 (except no toluene in sample 011). Significant benzene was also detected in sample 013. Styrene (14J µg/kg) was detected in sample 012.

As was the case with the surface soil samples, the cyanide contamination can be attributed to the FMGP. It is probable that the deep soil PAH contamination came from both the FMGP and the petroleum tank farm. The presence of mixed BTEX compounds and PAHs indicates a mixed source, as volatiles are not considered abundant in coal tar.

7.3 GROUND WATER SAMPLES

Ground water sampling data indicated no significant dissolved metals contamination when compared to background concentrations. Some organics were detected; however, in sample 204, piezometer well located downgradient of the site. This sample contained 3 PAHs: acenapthalene (65 µg/L), fluorene (25 µg/L), and phenanthene (46 µg/L); and 93 µg/L benzene (Figure 6-2; Appendix C). Additionally, syanide was detected in all ground water samples at concentrations above the 17 µg/L detection limit: 520J µg/L in sample 201, 27J µg/L in sample 202, 590J µg/L in sample 203, and 1,600J µg/L in sample 204. Cyanide was not detected in the background sample.

7.4 SURFACE WATER AND SEDIMENT SAMPLES

Surface water and sediment was sampled at three locations downgradient of the site. Additionally, the East St. Louis raw water intake, located across the Mississippi River from the site was sampled

Table 7-2 PAHs and Cyanide in Subsurface Soil Samples and Product Sample Laclede Coal Gas Site St. Louis, Missouri E & E/FIT; Harch 1991 Sample Series DSX44 (mg/kg)

Coapound	(2-7') 001*	(7-11') 002*		(17-21') 004*				(8-12') 008	(4-6') 009	(12-14') 010	(11-17') 011		(11-12.5') 013	501
Haphthalene		0 . 40U				23	87	2200	68	0.89	30		55	700
2-methylnaphthalene		0.40U				80	43	69	84	0.98	23			4000
2-chloronapththalene		0.40U								***				450
Acenaphthylene		0.40U						15			20			450
Acenaphtene		0.40U						32	12	0.15	15		22	620
Fluorene		0.40U				20	'	J	13	0.17	24		21	830
Phenanthrene	1.4	0.40U				63	20	J	46	0.51	71		59	. 300
Anthracene	0.38	0.40U	•~					J		-	23		18	110
Fluoranthene	2.0	0.52	0.54	-	_						180		39	620
Pyrene	2.8	0.53	0.80			48	J	J	J	J	J		65	2800
Benzo(a)anthracene	1.8	0.40U				18		J			99		21	1600
Chrysene	1.8	0.40U	Ö.44			18		J	16	0.18	94	_	21	2700
Benzo(b) fluoranthene	1.6	0.400						61			76		15	_
Benzo(k)fluoranthene	1.4	0.40U						52			89		17	
Benzo(a)pyrene	1.5	0.40U						95			120		23	970
Indeno(1,2,3-CD) pyrene	0.93	0.40U						37			75			_
Dibenso(a,h)anthra- cene		0.40U								-				
Dibenso(g,h,i)pery- lene	1.2	0.40U			<u></u>			44			88			
Cotal PAH Content	16.8	1.1	1.8			270	150 .	2605	239	2.9	1027		376 20	0,850
Cyanide Hassassassassassassassas	26	24				87	180	860		48		170		

^{* =} Background sample

Mote: See Figure 6-2 for sample locations and Appendix D for complete analytical results.

^{- =} Undetected (U) with detection limit given for background sample 002.

J = Detected, but results are invalid.

(304 and 304D). The most upgradient of these samples (302 and 402) generally exhibited the most contamination, though concentrations are fairly comparable (Tables 7-3 and 7-4). An exception is total chromium (12-14 µg/L), lead (15-18 µg/L), and zinc (46-54 µg/L) detected in samples 304 and 304D, collected from the surface water intake. These metals were not detected in the dissolved fraction of these samples. The only organics detected in these samples are relatively low levels of PAHs and cyanide found in the sediment samples (Table 7-4). A background sediment and surface water sample was not collected, so that no comparisonn to background can be made for these metals. However, sample concentrations are relatively low as compared to other samples collected.

7.5 PRODUCT SAMPLE

The oil sample (501) collected from an 80,000-gallon tank (Figure 2-3, tank #7) contained significant concentrations of many PAHs (Table 7-2). Attribution of PAH contamination on site is very difficult to assess, based on the fact that in 1986, PF & T experienced a spill of the same oil (#6) which was sampled, over the entire site. It was also common practice for FMGPs to bury their tar wastes on site. PAHs are a major constitudent of coal tar wastes. Thus, surface and subsurface soil PAH contamination could be attributable to either PF & T activities or the FMGP.

Table 7-3
Selected Total Metals
Sediment and Surface Water Samples
Laclede Coal Gas Site
St. Louis, Missouri
E & E/FIT; March 1991
Sample Series DSX44

Sample #	As	Ba	Cu	Cr	Ni	Pb	Se	V	Zn		
	Sediment Samples (mg/kg)										
401	3.7	140	9.1	8.7	10	30J	,	15	35J		
401D	4.0	140	8.2	9.0	11	13J	,	17	36J		
402	8.4	160	26	16	18	36J	2.0J	27	77J		
403	7.1	160	23	12	16	31J	,	25	64J		
			Vate	r Sampl	es (µg/l	.)					
301	,	,		,		7.0	,				
301D	,		,	,		7.2					
302		,	,			9.7					
303		280				24U	I*	6	2 89บ		
304			,	14		18	,		54		
304D		,	,	12		15					

 $[\]star$ = Invalid analysis, but 11 µg/L dissolved selenium was detected in this sample.

Note: See Figure 6-2 for sample locations and Appendix D for complete analytical results.

U = Undetected; detection limit given if above detected values.

J = Results are reported, but invalid by approved QC procedures.

Table 7-4
PAHs, Cyanide, and Total Hydrocarbons
Sediment Samples
Laclede Coal Gas Site
St. Louis, Missouri
E & E/FIT; March 1991
Sample Series DSX44

		2535555555555555	***********	*********
(µg/kg)	401	401D	402	403
Phenanthene			2900	4400
Fluoranthene		750	5000	5100
Pyrene		960	8000	6400J
Benzo(a)anthracene		460	3500	4200
Di-n-octyl phthalate		470	3900	4900
Benzo(k)fluoranthene			2900	3100
Benzo(a)pyrene		430	2600	5600
Benzo(GHI)perylene		~	3500	4100
Cyanide			1600	
Total hydrocarbons	3100	3100U	8200	4900

J = Results are reported, but invalid by approved QC procedures.

Note: See Figure 6-2 for sample locations and Appendix D for complete analytical results.

U = Undetected, with detection limit given.

⁻⁻⁻ Indicated undetected with detection limits below detected values.

SECTION 8: SUMMARY AND CONCLUSIONS

A screening site inspection was conducted at the Laclede Coal Gas site, St. Louis, Missouri. The Laclede site is a FMGP presently occupied partially by a petroleum tank farm. The site was originally investigated as the Mound Street Power Plant, due to observed oily waste releases to the adjacent Mississippi River. Previous oil sampling has not detected polychlorinated biphenyls (PCBs). The Laclede FMGP is thought to be the largest FMGP in EPA Region 7. The presence of a surface water intake located almost directly across the Mississippi River from the site. prompted EPA to request an SSI of the site. The purpose of the investigation was to evaluate the potential threat this site poses to the ground water, surface water, air and on-site pathways. These objectives were addressed through both non-sampling and sampling data collection strategies. Non-sampling data allowed the development of a detailed background on all potential targets. The sampling allowed identification of on-site contaminants, and gathered data on the areal and vertical extent of contamination.

Non-sampling data indicated that the only potential primary targets for the site included a population of approximately 300,000 people who receive drinking water from the intake located directly across the Mississippi River from the site. Sampling data, however, indicated that these people are secondary targets because no significant contamination was found in the surface water samples (including an intake sample).

Thirteen subsurface soil, twelve surface soil, six surface water, four sediment, six ground water, and one oil sample were collected and submitted to the CLP for volatiles, semi-volatiles, metals, and cyanides analysis. Additionally, the FASP screened 92 samples on-site to help select CLP sample locations. Seventeen drums of IDW were generated and a subcontractor used for soil sampling.

The sampling data identified cyanides, BTEX compounds, and PAHs as the major on-site contaminants. The cyanide contamination was found throughout the site in the 0 to 2 feet deep soil samples. Deeper soil samples revealed cyanide contamination at depths as great as at least 11 feet. PAH contamination, both in shallow and deep soil samplings, was restricted to small areas within the site as evidenced by FASP data and

was found at depths as great as 38 feet. The greatest area of PAH contamination was found within the bermed tank farm (Figure 6-2). Only one ground water sample (204) showed PAH contamination, but all contained cyanide. Benzene was also detected in sample 204.

Attribution is difficult to assess because PF & T currently stores #6 oil on site in large quantities. The total PAH content of this oil is very high. PF & T experienced a spill of this oil over the entire site in 1986. BTEX compounds were found at relatively high concentrations in several subsurface soil samples, indicating BTEX wastes are sometimes mixed with PAH wastes. It is known tht FMGPs frequently buried coal tar waste on site. PAHs are a major constituent of coal tars. Thus, both the FMGP and PF & T could be responsible for the PAH contamination found abundantly in on-site soils, and to a much lesser extent in the ground water. The cyanide contamination, however, is assumed to largely be a result of the FMGP as cyanide is not considered a significant constituent of most oils.

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APPENDIX A Technical Directive Document

P	Cost Contest &
H	Live Target State
3184	Apparel Miniberts
į.	

FIT ZONE II CONTRACT Contract Number 68-01-7347

	TOD Number.
F	097-9008-020

2A. Amendment:

	TECHNICA	IL DIRECTI	VE DOCUME	NI (IDD)		Administrative Technical
3A. Priority: High	3B. Key EPA Contact					
Medium ☐ Low	Name: GREG R	EESOR		Phone	:551-76	95
4A. Estimate of Technical Hours:	4B. Subcontract: T&M IDW DISPOSAL	4C. Estimate Subcont	of ract Cost:	5A. SSID Numbe	r: 58. CEF	RID Number:
*1315	T&M DRILLING	*NTE 35,	882 22,00	O UNASSIGNE	D MOD	81715980
5C. EPA Site Name:		ı	5D. City/Cou	nty/State:		
LACLEDE COAL GAS	SITE	·	ST. LOU	IS/ST. LOUIS/	MISSOURI	
	Standard Report Othe	er (Specify):	7A. Activity		7B. Estimated (Completion
Letter Report	Formal Briefing			8/29/90	10/31/	91
8A. Type of Activity:		nforcement Su	• •	raining	8B. FIT/SCAP Will Deliver a Unit of the	rable Meet
∑SI □ RCRA-SI □ ESI	, _	rogram Manage quipment Main		eneral Technical ssistance	▼ Yes	⊒No
	tion: PERFORM A SCRE	ENING SITE	E INVESTIGA	TION OF THE 1	LACLEDE COAL	GAS SITE
LOCATED IN ST. L					WED.	
					NAFILL	
				N a.)	
				SEP 1	E 1991	
10. Specific Elements:	1. REVIEW BACKGR	OUND INFO	RMATION	FRE	(11) nuirim De	adlines:
2. PREPARE DRAF	T HRS SCORE	·····				
3. CONDUCT SITE	RECONNAISSANCE		·		<u> </u>	
4. PREPARE WORK	PLAN			···		
5. PREPARE BID	SPECS AND SUBCONTE	ACT PACKA	GES			
6. IMPLEMENT WO	RK PLAN AND PRODUC	E TRIP AND	D FINAL RE	PORTS (DRAFT)	6. EST	10/15/91
7 *OVERSEE SUBC	CONTRACTOR SAMPLE A	ND DISPOS				
		<u> </u>		al Scope Attached	<u> </u>	
_	NDED TO ADD TWO SI		•		S_TO_SITE_(6	O HOURS)
UNLY DEAFT FINAL	REPORT IS EXPECTE	ED BY CONT	KACT_COMPL	ETION.		
· · · · · · · · · · · · · · · · · · ·						
13. Authorizing:	A6 ()		ARPO		14. Date:	
	(Signature)		—— □PO)	9/161	9/
THE SAME	发展 罗黎		ARAcc		16, Date:	
Maria Series	was the Mi			epted with eptions (Attached)	9/197	劳 经越
次次发展的影响	onunction FITOM Signature	门数字传播	PARTITION	octed - The second	THE PERSON NAMED IN	通过可能

APPENDIX B EPA Site Inspection Form 2070-13

		POTENTIAL	L HAZAKDOUS	WASTE	SITE		I.	1. IDENT			1
EPA		SITE	INSPECTION	REPOR	•		[·	01 STATE	02 SIT	E NUMBER	
	PA.s.	1 - SITE LOC	ATION AND II	NSPECT:	ON IN	ORMATION	L				
II. SITE NAME AND											
1 SITE NAME (Lega Laclede Coal Gas		scriptive name	e of site)	l.		ROUTE NO., C		CIFIC LOC	MOITA	IDENTIFIE	SR .
13 CT#U				04 ==	TE A	ZID CONTIN	6 600		Τ.	7	00 0000
)3 CITY				MO	- 1	ZIP CODE 06	o count	4.2	١٥	7 COUNTY CODE	DIST
St. Louis O COORDINATES		10 TYPE OF OW	NEBSUID (Ch.			.3103					
LATITUDE	LONGITUDE	i			•	r e	TATE	D CO!-	NTV	& Mare	ICTEST
į.	090° 11' 00.0"	F. OTHER						B. COU	_	E. MUN	CIFAL
111. INSPECTION 1		· OIRER						G. UNK			
1 DATE OF INSPECT		02 SITE STATU	S 03 Y	EARS O	POPER	MOITA					
3/3/91		_x_ACTIV		888		1 1945		<u>-</u>	UNK	NOWN	
MO/DAY/YR		INACT	l	EGINNI	NG YEA	· —	NG YEA	JR .	— ·		}
	NG THEODOG	Charle all Attent	Apples 1								
A FRA - B			••••	т.	D 1400	TCTD11	D 1 ^				Į
A. EPA XB.	EPA CONTRACTOR			-	J. MUN	LCIPAL CONT	~~CTOR		· -	190	<u> </u>
r emimo .	r. etime commi	(Name of fir	es j	_	Udna.					(Name of	rirm)
L. STATEI	. STATE CONTRAC		f firm)	G.	-THER		10	of for 1			1
05 CHIEF INSPECTO	 -		f firm) 06 TITLE		···	07 ORGANIZ	(Spec	Y1	In-	TELEPHO	NE NO
Keith A. Brown	. .	Į.	Environme	ntal		1	E/FIT		- 1	5 ТЕЦЕРНО (913) 432	
A. DIOWN			Scientist								-2501
09 OTHER INSPECTO	RS		10 TITLE			11 ORGANIZ	MOITA		12	2 TELEPHO	NE NO.
Anne Melia			Chemist			E &	E/FIT			(913) 432	-9961
Chris Williams	 		Geologist	:		E &	E/FIT			(913) 432	-9961
Well McCall			Geologist	:		E &	E/FIT			(913) 432	-9961
John Peck John Strobel			Geologist Chemist			E 4	E/FIT			(913) 432	-9961
Patty Roberts			Environme Scientist	ntal		E &	E/FIT			(913) 432	-9961
13 SITE REPRESENT.	ATIVES INTERVIEW	VED	14 TITLE			15 ADDRESS	5		1	6 TELEPHO	NE NO.
Randel Lewis			Terminal	Manage	r	Foot of	Mullar	nphy Stree	et	(314) 621	-0522
										()	
										()	
										()	
										()	
										()	
										()	
17 ACCESS GAINED (Check one)	BY 18 TIME OF	INSPECTION	19 WEATHER	CONDI	CIONS	· 					
x PERMISSION	. NA										
WARRANT											
			<u> </u>								
IV. INFORMATION	AVAILABLE FROM		103						- 1	3 ====	
01 CONTACT			02 OF (Age)			Lion)			10	3 TELEPHO	
Greg Reesor			EPA/S	Superf	ari Q					(913) 55	1-/695
04 PERSON RESPONS	SIBLE FOR SITE I	NSPECTION FORM	05 AGENCY	0	ORGA	NIZATION	07 TE	LEPHONE N	0. 0	8 DATE	
Keith A. Brown			E & E]	1	FIT	(913) 432-996	i	10/30/91	

EPA FORM 2070-13 (7-81)

EPA			SITE INSPECT	COM	REPORT			01 STATE		2 SITE NUMBER
		* 4	PART 2 - WASTE				ł	MO		2 SITE NUMBER D981715980
II. WASTE STA	TES, QUANTITIES, A									
01 PHYSICAL S	TATES	02 WAS	TE QUANTITY AT SI	TE	03 WASTE CHARACT	TERISTICS				
(Check all th	at apply)	(Measu	res of waste quant oust be independent	<u>: 1 - </u>	(Check all that	apply)				
xA. SOLID	E. SLURRY	C105	ust be independent	ا ''	x A. TOXIC	E. S	OLUBLE	3	г. н	IGHLY VOLATILE
B. POWDER,	FINES x F. LIQUID	Ì		Ì	B. CORROSIVE	. <u> </u>	NFECTIO	ous —	J. E	XPLOSIVE
x c. SLUDGE	G. GAS	l	TONS Unknown	- I	C. RADIOACT	IVE G. F	LAMMABI	LE I	K. R	EACTIVE
D. OTHER		CUBIC	YARDS		x D. PERSISTER	чт н. 1	GNITABI	LE I	L. I	NCOMPATIBLE
_ 	(Specify)			_					A. N	OT APPLICABLE
		NO. OF	DRUMS	-				_		
III. WASTE TO	PE								<u> </u>	
CATEGORY	SUBSTANCE NAME		01 GROSS AMOUNT	02	UNIT OF MEASURE	E 03 COMM	ENTS			
SLU	SLUDGE	*	Unknown	\prod		Coal ta	rs			
OTM	OILY WASTE									
SOL	SOLVENTS									
PSD	PESTICIDES									
occ	OTHER ORGANIC CHEM	ICALS		Г						
IOC	INORGANIC CHEMICAL	s	Unknown			Purific	r wast	0.8		
ACD	ACIDS			T						
BAS	BASES			Т						
MES	HEAVY METALS	-		Г						
IV. HAZARDOUS	SUBSTANCES (See A	ppendi	k for most frequen	tly	cited CAS Numbe	rs)				
01 CATEGORY	02 SUBSTANCE NAM	E	03 CAS NUMBER	0.	STORAGE/DISPOS	AL METHOD	05 CON	CENTRATI	ON	06 MEASURE OF CONCENTRATION
IOC	Cyanides	·	143-33-9	B	rial		>500			ppm
			151-50-8	L						
	<u></u>			丄			<u> </u>			
SLU	Poly Aromatic		8007452	B	urial		>100			bbw
	hydrocarbons (PA	Hs)		丄			<u> </u>			
l 				\perp						
				丄				 		<u> </u>
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<u></u>				↓_			L			
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				┷			<u> </u>			<u> </u>
	<u> </u>		<u> </u>	丄			<u>. </u>			<u> </u>
	S (See Appendix for			 -				·,		
CATEGORY	01 FEEDSTOCK	NAME	02 CAS VIMBER		CATEGORY	01 FEED	STOCK N	IAME	°	2 CAS NUMBER
FDS					FDS				_	
FDS			}		PDS					
FDS	_				FDS	·				
FDS					FDS	_			<u> </u>	
VI. SOURCE	S OF INFORMATION (ite sp	ecific references	, e.	g., state files.	, sample an	alysis,	reports	<u> </u>	
E & E/FIT	files									
	;									
1										

EPA FORM 2070-13 (7-81)

	POTENTIAL HA	REARC	DOUS WASTE SITE		TIPICATION	
EPA	SITE INS	SPECT	TION REPORT	01 STATE	02 SITE NU D981715	MBER
	PART 3 - DESCRIPTION OF	HAZI	ARDOUS CONDITIONS AND INCIDENTS	L	1	
II. HALARDOUS CONDITIONS	AND INCIDENTS					
01 x A. GROUND WATER CONT	ramination	02	X OBSERVED (DATE: 3/1991)	x POT	ENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED: None known	04 8	NARRATIVE DESCRIPTION			
Three PAHs and benzene with the known to be used for dri	were detected in a downgra inking.	adier	nt well. No ground water in the	surround	ing area is	
01 x B. SURFACE WATER CON	NTAMINATION	02	OBSERVED (DATE:)	x POT	ENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED: 300,000	04 1	NARRATIVE DESCRIPTION			
There is a strong potent was detected immediately across the Mississippi l	tial for a ground water to y downgradient of the site River.	0 SUI 0.]	rface water release of contamina Low potential for release to sur	tion, how face wate	ever, no co r intake lo	ntsmination cated
01 x C. CONTAMINATION OF	AIR	02	OBSERVED (DATE:)	x POT	ENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED:	04 1	NARRATIVE DESCRIPTION			-
Low potential for air co	ontamination due to site's	a gr	avei covering.			
01D. FIRE/EXPLOSIVE C	ONDITIONS	02	OBSERVED (DATE:)	PO1	ENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED:	04	NARRATIVE DESCRIPTION			
None known in relation Ol x E. DIRECT CONTACT	CO FRISE.	02	x OBSERVED (DATE: 3/1991	<u>x</u> P01	CENTIAL	ALLEGED
	AFFECTED: on-site worker: e site boundaries are con-		04 NARRATIVE DESCRIPTION nated with cyanide and PAHs.			
01 x F. CONTAMINATION OF	SOIL	02	x OBSERVED (DATE: 3/1991	PO'	TENTIAL	ALLEGED
03 AREA POTENTIALLY AFFEC	· · · · · · · · · · · · · · · · · · ·	04	NARRATIVE DESCRIPTION			
Cyanide in the 10 to 10	(Acres) Os ppm range. Total PAHs	in	the 10 to 1000s ppm range.			
01 G. DRINKING WATER C		02	OBSERVED (DATE:	PO	TENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED:	04	NARRATIVE DESCRIPTION	_		•
None known.						
01 x H. WORKER EXPOSURE/	INJURY	02	OBSERVED (DATE:) PO'	PENTIAL	ALLEGED
03 WORKERS POTENTIALLY AF	FECTED:		NARRATIVE DESCRIPTION			_
See Direct Contact abou	<u></u>					
01 I. POPULATION EXPOS	SURE/INJURY	02	OBSERVED (DATE:) PO	TENTIAL	ALLEGED
03 POPULATION POTENTIALLY	AFFECTED:		NARRATIVE DESCRIPTION			-
None known.						
I						

	AL HAZAKDOUS WASTE SITE	1. IDENTIFICATION
EPA SIT	E INSPECTION REPORT .	01 STATE 02 SITE NUMBER
PART 3 - DESCRIPTION	N OF HAZARDOUS CONDITIONS AND INCL	DENTS MO D981715980
II. HAZARDOUS CONDITIONS AND INCIDENTS (CONT	INUED)	
01 J. DAMAGE TO FLORA	02 OBSERVED (DATE:) POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION		
None observed.		
101.6 ODB011-0-1.		
01K. DAMAGE TO FAUNA	02OBSERVED (DATE:) POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION (Include name(s) of s	pecies)	•
None observed.		
		•
A4 COMPLUTIVETOR OF COAR CULTU	03	
01 x L. CONTAMINATION OF FOOD CHAIN	02 OBSERVED (DATE:) x POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION		
This may occur if surface water is contamina water contamination.	ited, however, present sampling did	not detect significant surface
01 x M. UNSTABLE CONTAINMENT OF WASTES	02 OBSERVED (DATE:) x POTENTIAL ALLEGED
(Spills/runoff/standing liquids/leaking drum		
03 POPULATION POTENTIALLY AFFECTED:	04 NARRATIVE DESCRIPTION	
Waste may be land applied or build in unline		
01 N. DAMAGE TO OFF-SITE PROPERTY	02 OBSERVED (DATE:) POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION		
None known.		
-		
At a salmiury men of street	03	
01 O. CONTAMINATION OF SEWERS, STORM DRAINS, WWTPB	02 OBSERVED (DATE:	POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION		
None known.		
01P. ILLEGAL/UNAUTHORIZED DUMPING	02 OBSERVED (DATE:) POTENTIAL ALLEGED
04 NARRATIVE DESCRIPTION		
None known.		
1 :		
05 DESCRIPTION OF ANY OTHER KNOWN, POTENTIAL,	OR ALLEGED HAZARDS	
		is an shandoned numb bouse
E & E/FIT observed oil and coal tars leach; pipes. The pipes have been sealed with con	crete, but contaminants are leachi	ng through the concrete plugs.
III. TOTAL POPULATION POTENTIALLY AFFECTED:	300,000	
TW COMMENTS		
IV. COMMENTS		·
1		
V. SOURCES OF INFORMATION (Cite specific ref	erences, e.g., state files, sampl	e analysis, reports)
E & E/FIT on site reconnaissance, November E & E/FIT files	20, 1990	
E & E/PIT files E & E/PIT SSI, March 3-9, 1991		

	POTENTIAL HALI	ARDOUS W	aste site		I. IDE	ENTIFICATION
EPA	SITE INS.	LTION N	EPORT.		01 STAT	2 02 SITE NUMBER D981715980
	PART 4 - PERMIT AND	DESCRIP	TIVE INFO	RMATION	MO	D981715980
II. PERMIT INTORMATION						
01 TYPE OF PERMIT ISSUED	02 PERMIT NUMBER	03 DATE	ISSUED	04 EXPIRATION DATE	105 CC	MMENTS
(Check all that apply)	· · · · · · · · · · · · · · · · · · ·	l				,
- -		{	İ			}
A. MPDES	i	1			1	
B. UIC						
C. AIR						
D. RCRA						
E. RCRA INTERIM STATUS						
F. SPCC PLAN						
G. STATE (Specify)						
H. LOCAL (Specify)						
I. OTHER (Specify)		<u> </u>				
x J. NONE KNOWN				<u> </u>		
III. SITE DESCRIPTION						
01 STORAGE/DISPOSAL 0	2 AMOUNT 03 UNIT OF	MEASURE	04 TREAT	Ment		05 Other
(Check all that apply)			(Chec	k all that apply)	- 1	
A. SURFACE IMPOUNDMENT			λ. I	ncineration	- 1	
B. PILES				NDERGROUND INJECTI	ON I	x A. BUILDINGS ON SITE
C. DRUMS, ABOVE GROUND		_		HEMICAL/PHYSICAL	···	
	Togget land use			•	Į.	
	resent land use		_	IOLOGICAL	_	06 AREA OF SITE
E. TANK, BELOW GROUND			E. W	ASTE OIL PROCESSIN	G [
F. LANDFILL			F. S	OLVENT RECOVERY	i	~7 (Acres)
g. Landfarm			a. o	THER RECYCLING/REC	OVERY	i
H. OPEN DUMP			н. о	Ther	[
x I. OTHER Buried pits Unk	nown FMGP			(Specify)		
(Specify)					1	
107 COMMENTS						
A petroleum tank farm curren plant rubble and other const	tly occupies the site ruction debris underli	Subsui es the	rface exp site.	loration reveals t	hat mos	t of the old coal gas
IV. CONTAINMENT						
1 CONTAINMENT OF WASTES (Chec	rk one)					
A. ADEQUATE, SECURE	B. MODERATE	c. I	NADEQUATE	, POOR _x_D. I	NSECURE	, unsound, dangerous
42 2222222						
102 DESCRIPTION T DRUMS, DIKIS Coal tar and purifier waste			site or s	old.		
1						i
				-		
V. ACCESSIBILITY						
02 COMMENTS	ABS _x NO					
Most of the site is covered	with gravel.					
VI. SOURCES OF INFORMATION (cite specific reference	es, e.g.	, state i	iles, sample analy	ysis, r	eports)
E & E/FIT files E & E/FIT SSI, March 3-9,	1991				-	
3 - 2/12 - 202, 100 - 30 - 37,						
}						
1						
1						

	POTENTIAL HAZARDOU	S WASTE SIT	B		I. IDENT	IFICATION	
EPA	SITE INSPECTIO	N REPOR			01 STATE	02 SITE NUMBER D981715980	
	ATER, DEMOGRAPHIC,	AND ENVIRO	MENTAL D	ATA		D981715980	
II. DRINKING WATER SUPPLY							
01 TYPE OF DRINKING SUPPLY (Check as applicable)	02 STATUS	NA			03 DISTANC	han nine miles	up-
SURFACE WELL COMMUNITY A. x B.	A	B	MONITO	-	Mississip A.	E directly acr pi River. (mi)]
NON-COMMUNITY C. x D	D	E	F	-	В. —	(ini)	'
III. GROUND WATER						· · · · · · · · · · · · · · · · · · ·	
01 GROUND WATER USE IN VICINITY (Check	one)						
COMMERCIAL,	es available) INDUSTRIAL, IRRIGA ter sources availa	(Limit TION	ERCIAL, 1 ed other	NDUSTRIAI BOUICOB 8	L IRRIGATIO Ivailable)		OT USED,
02 POPULATION SERVED BY GROUND WATER N	one known	03 DISTANCE	TO NEARI	ST DRINK	ING WATER W	ELL Unknown	(mi)
04 DEPTH TO GROUNDWATER 05 DIRECTION OF 25-30 (ft) Southeast	GROUNDWATER FLOW	06 DEPTH TO OF CONCE 25		07 POTENT OF AQU Unknown	TIAL YIELD UIFER (gpd)		X NO
09 DESCRIPTION OF WELLS (Including usag No wells are known within four miles surface water.	· · ·					•	ndance of
10 RECHARGE AREA		11 DISCHARG	E AREA	 			
YES COMMENTS		_x_YES	COMMENT	5			ŀ
x NO		мо					
IV. SURFACE WATER							
01 SURFACE WATER USE (Check one)		·					
* A. RESERVOIR, RECREATION B. IRE DRINKING WATER SOURCE IMP	RIGATION, ECONOMICA PORTANT RESOURCES	LLY x C.	COMMERCI.	AL, INDUS	TRIAL	. NOT CURRENT	LY USED
02 AFFECTED/POTENTIALLY AFFECTED BODIES	OF WATER						
NAME:			λ	FFECTED	מ	STANCE TO SIT	E
Mississippi River					< 1/	<u> </u>	(mi)
					-		—— (mi)
				_			(mi)
V. DEMOGRAPHIC AND PROPERTY INFORMATIO	ON						
01 TOTAL POPULATION WITHIN			02	DISTANCE	TO NEAREST	POPULATION	
ONE (1) MILE OF SITE TWO (2) MILES	OF SITE THREE (3) MILES OF	SITE	<	1/4	(mi)	
A. 6586 B. 32,564	c. <u>100,</u>	924	_				•
NO. OF PERSONS NO. OF PER		. OF PERSON	s	· · · · · · · · · · · · · · · · · · ·		<u> </u>	
03 NUMBER OF BUILDINGS WITHIN TWO (2)	AILES OF SITE	04 DISTA	NCE TO NE	AREST OFF	-SITE BUIL	DING	
Heavy industrial area	-		< 1/2		(mi)		
05 POPULATION WITHIN VICITY OF SITE	(Provide narrative e.g., rural, vill	description	n of natu y populat	re of por ed urban	ulation wi	thin vicinity	of site,
Densely populated urban area.							
,							

EPA FORM 2070-13 (7-81).

	POTENTIAL HAZARDO	OUS WASTE SITE		I. IDENTIF	
ВРА	SITE INSPECTI	· · · · · · · · · · · · · · · · · · ·	•	01 STATE 02	SITE NUMBER D981715980
	WPIER, DEMOGRAPHIC	, AND ENVIRON	MENTAL DATA		D34272330V
VI. ENVIRONMENTAL INFORMATION					
01 PERMEABILITY OF UNSATURATED ZONE (Theck one) Un	known			
A. 10 ⁻⁶ - 10 ⁻⁸ cm/secB. 10 ⁻	-4 - 10 ⁻⁶ cm/sec	c. 10 ⁻⁴ -	10 ⁻³ cm/sec	D. GREATER T	THAN 10 ⁻³ cm/sec
02 PERMEABILITY OF BEDROCK (Check one					
A. IMPERMEABLEB. RELA:	TIVELY IMPERMEABLE	C. RE	LATIVELY PERMEABLE	E	. VERY PERMEABLE
(Less than 10 ⁻⁶ cm/sec) (10 ⁻⁴			² - 10 ⁻⁴ cm/sec)	(Great	er than 10 ⁻² cm/sec)
	NTAMINATED SOIL ZOI	NE 05	SOIL PH		•
	(ft)		nknown		· · · · · · · · · · · · · · · · · ·
06 NET PRECIPITATION 07 ONE YEAR 24	HOUT RAINFALL 0	SITE SLOPE	DIRECTION OF SIT	E SLOPE TER	RRAIN AVERAGE SLOPE
0 (in) 3	(in)	0~5	Easterly		1-2
09 FLOOD POTENTIAL	10 STEE TS (ON BARRIER TET	AND COLUMN HIGH	W17170 1001	A, RIVERINE FLOODWAY
SITE IS IN 500 YEAR FLOODPL	AIN	ON BARRIER ISE	AND, COASTAL HIGH	HAZARD AKEA	A, RIVERINE PLOODWAY
11 DISTANCE TO WETLANDS (5 acre minim	nw)	12 DISTANCE T	O CRITICAL HABITA	T (of endang	gered species)
	other			< 1	(mi)
A. greater than 4 (mi) B. g	reter than 4 (mi)	Endanger	ED SPECIES: Sick	lefin Chub	
13 LAND USE IN VICINITY					
DISTANCE TO: RESIDEN	TIAL AREAS; NATION. RESTS, OR WILDLIFE	AL/STATE PARKS	s. a	GRICULTURAL	LANDS AG LAND
COMMERCIAL/INDUSTRIAL FO	RESTS, OR WILDLIFE	RESERVES			
A. <1 (mi)	B. < 1	(mi)	с	Unknown (mi) D	(mi)
4 DESCRIPTION OF SITE IN RELATION TO	SURROUNDING TOPOG	RAPHY			
	# Abs Missississ	.			
Site is located on the flood plain o	r the mississippi	Kiver.			
·	-				
					
VII. SOURCES OF INFORMATION (Cite sp	ecific references,	e.g., state	files, sample anal	ysis, repor	ts)
E & E/FIT files					
\					
<u> </u>					
1					

EPA FORM 2070-13 (7-81)

	PO	TENTIAL LAREDUS WASTE SITE	I. IDENTIFICATION
EPA		SITE INSPECTAON REPORT	01GE 02 SITE NUMBER D981715980
	PART	6 - SAMPITANO FIELD INFORMATION	
II. SAMPLES TAKEN			
SAMPLE TYPE	01 NUMBER OF SAMPLES TAKEN	02 SAMPLES SENT TO	03 ESTIMATED DATE RESULTS AVAILABLE
GROUND WATER	6	CLP	Current
SURFACE WATER	6	CLP	Current
WASTE			
AIR			
RUNOFF			
SPILL			
SOIL/Sediment	25/4	CLP	Current
VEGETATION			
OTHER (Product)	1	CLP	Current
III. FIELD MEASUREME	NTS TAKEN		
01 TYPE FASP	02 COMMENTS 92 soil, 3 seding a mobile lab.	ment, and 4 surface water samples were confidence to the confidence of the confidenc	llected and screened on site in analysis.
Parameters	pH, conductivity	y, and temperature were recorded for grou	and surface water samples.
1			
IV. PHOTOGRAPHS AND	Maps		
01 TYPE x GROUND	AERIAL	02 IN CUSTODY OF E & E/FIT	
	<u></u>	(Name of organization	on or individual)
03 MAPS 04	LOCATION OF MAPS		
	& E/FIT		
V. OTHER FIELD DATA	COLLECTED (PEGA109	narrative description)	
		•	
1		•	•
l			
1			
1			
	RMATION (Cite specif	ic references, e.g., state files, sample	analysis, reports)
E & E/FIT files	 -		
l			

		POTI	ENTIAL HARARDOUS	WASTE SILE	I. IDENTIF	ICATION
BPA			SITE INSPECTION	REPROT	01 STATE 02	SITE NUMBER D981715980
		P	ART 7 - OWNER IN			
II. CURRENT OWNER(S)				PARKET COMPANY (If applicat		
01 NAME Petroleum, Fuel & Terminal	с.	02 D	+B NUMBER	08 NAME Apex Oil		09 D+B NUMBER
03 STREET ADDRESS (P.O. BOX, Poot of Mullanphy	RFD #, ET	c.)	04 SIC CODE	10 STREET ADDRESS (P.O. BOX	, RFD #, ETC	.) 11 SIC CODE
05 CITY	06 STATE	07 Z	IP CODE	12 CITY	13 STATE	14 ZIP CODE
St. Louis	мо	63:	105	į		
01 NAME		02 D	+B NUMBER	08 NAME		09 D+B NUMBER
03 STREET ADDRESS (P.O. BOX,	RFD 0, ET	(C.)	04 SIC CODE	10 STREET ADDRESS (P.O. BOX	, RFD #, ETC	:.) 11 SIC CODE
05 CITY	06 STATE	07 Z	IP CODE	12 CITY	13 STATE	14 ZIP CODE
)1 NAME		02 D	+B NUMBER	08 NAME		09 D+B NUMBER
)3 STREET ADDRESS (P.O. BOX,	RFD #, E	rc.)	04 SIC CODE	10 STREET ADDRESS (P.O. BOX	, RFD #, ETC	c.) 11 SIC CODE
05 CITY	06 STATE	07 Z	IP CODE	12 CITY	13 STATE	14 ZIP CODE
III. PREVIOUS OWNER(S) (List	most re	cent	first)	IV. REALTY OWNER(S) (If app	licable; li	st most recent fix
01 NAME Laclede Gas & Light Co.		02	D+B NUMBER	01 NAME	· · · · · · · · · · · · · · · · · · ·	02 D+8 NUMBER
u3 STREET ADDRESS (P.O. Box, Unknown	RFD ♦, ●	tc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box	, RFD #, et	c.) 04 SIC CODE
5 CITY	06 STAT	E 07	ZIP CODE	05 CITY	06 STAT	E 07 ZIP CODE
1 NAME Mound Street Warehouse Cor	<u>.</u> P.	02	D+B NUMBER	01 HAME		02 D+B NUMBER
03 STREET ADDRESS (P.O. Box, Unknown	RFD #, •	tc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box	r, RFD #, et	e.) 04 SIC CODE
US CITY	06 STAT	E 07	ZIP CODE	05 CITY	06 STAT	E 07 ZIP CODE
1 NAME		02	D+B NUMBER	01 NAME		02 D+B NUMBER
"3 STREET ADDRESS (P.O. Box,	RFD #,	tc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box	k, RFD #, et	c.) 04 SIC CODE
05 CITY	06 STAT	E 07	ZIP CODE	05 CITY	06 STAT	E 07 ZIP CODE
7. SOURCES OF INFORMATION (Cite spec	ific	references, e.g	., state files, sample analy	sis, reports	1)
E & E/FIT files				·		

	POTENTIAL HAZARDOUS		I. IDENTIFIC	
EPA	SITE INSPECTION	REPORT	01 STATE 02 S	ITE NUMBER
	PART 8 - OPERATOR IN	FORMATION	HO I	3961713960
II. CURRENT OPERATOR (Provide if dif	ferent from owner)	OPERATOR'S PARENT COMPANY	If applicable)
01 NAME	02 D+B NUMBER	10 NAME	} :	11 D+B NUMBER
NA				
03 STREET ADDRESS (P.O. BOX, RFD #,	ETC.) 04 SIC CODE	12 STREET ADDRESS (P.O. BO)	, RFD #, ETC	.) 13 SIC CODE
05 CITY 06 STATE	07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION 09 NAME OF	OWNER		· ·	
III. PREVIOUS OPERATOR(S) (List most vide only if	recent first; pro- different from owner)	PREVIOUS OPERATORS' PARENT	COMPANIES (I	f applicable)
01 NAME	02 D+B NUMBER	10 NAME		11 D+B NUMBER
03 STREET ADDRESS (P.O. Box, RFD #,	etc.) 04 SIC CODE	12 STREET ADDRESS (P.O. Bo	x, RFD #, etc	.) 13 SIC CODE
05 CITY 06 STAT	E 07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION 09 NAME OF OWN 1888-1945	ER DURING THIS PERIOD		1	
01 NAME	02 D+B NUMBER	10 NAME		11 D+B NUMBER
03 STREET ADDRESS (P.O. Box, RFD #,	etc.) 04 SIC CODE	12 STREET ADDRESS (P.O. Bo	x, RFD #, etc	:.) 13 SIC CODE
05 CITY 06 STAT	E 07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION 09 NAME OF OWN	ER DURING THIS PERIOD			
01 NAME	02 D+B NUMBER	10 NAME		11 D+B NUMBER
03 STREET ADDRESS (P.O. Box, RFD #,	etc.) 04 SIC CODE	12 STREET ADDRESS (P.O. Bo	x, RFD *, etc	:.) 13 SIC CODE
05 CITY 06 STATE	TE 07 ZIP CODE	14 CITY	15 STATE	16 ZIP CODE
08 YEARS OF OPERATION 09 NAME OF OWN	VER DURING THIS PERIOD			
IV. SOURCES OF INFORMATION (Cite spe	cific references, e.g	., state files, sample analy	sis, reports)
E & E/FIT files				

		POT	ENT 1	AL HAZARDOUS	WASTE SITE		I. 16PWIFE		
EPA			SIT	E INSPECTION	REPORT .	Ţ	01 STATE 02	SITE	NUMBER 715980
	PART	9 –	GEN	erator/Transi	PORTER INFORMATION	L			
II. ON-SITE GENERATOR							<u>_</u> .		
01 NAME	ł	02 D	+B N	UMBER					- [
NA									į
03 STREET ADDRESS (P.O. BOX, 1	RFD #, ET	C.)	04	SIC CODE					
05 CITY	06 STATE	07	ZIP	CODE					
III. OFF-SITE GENERATOR(S)		<u> </u>		 -					
01 NAME NA		02	D+B	NUMBER	01 NAME	-		02 D	+B NUMBER
03 STREET ADDRESS (P.O. Box,	RFD #, et	:c.)	04	SIC CODE	03 STREET ADDRESS	(P.O. Box,	RFD #, etc.	,	04 SIC CODE
05 CITY	06 STATI	07	ZIP	CODE	05 CITY		06 STATE	07 Z	IP CODE
01 NAME	<u>1</u>	02	D+B	NUMBER	01 NAME			02 E	+B NUMBER
03 STREET ADDRESS (P.O. Box,	RFD #, e	tc.)	04	SIC CODE	03 STREET ADDRESS	(P.O. Box	, RFD #, etc.	.)	04 SIC CODE
05 CITY	06 STAT	E 07	ZIP	CODE	05 CITY		06 STATE	07 2	ZIP CODE
IV. TRANSPORTER(S)	L							L	
01 NAME NA		02	D+B	NUMBER	01 NAME			02 1	D+B NUMBER
03 STREET ADDRESS (P.O. Box,	RFD #, e	tc.)	04	SIC CODE	03 STREET ADDRESS	P.O. Box	, RFD #, etc	.)	04 SIC CODE
05 CITY	06 STAT	E 07	ZIP	CODE	05 CITY		06 STATE	07 :	ZIP CODE
01 NAME	<u> </u>	02	D+B	NUMBER	01 NAME			02 1	D+B NUMBER
03 STREET ADDRESS (P.O. Box,	RFD #, e	tc.)	04	SIC CODE	03 STREET ADDRESS	S (P.O. Box	, RFD #, etc	.)	04 SIC CODE
05 CITY	06 STAT	E 07	ZIP	CODE	05 CITY		re STATE	07	ZIP CODE
V. SOURCES OF INFORMATION (C	ite spec	ific	ref	erences, e.g.	, state files, sam	aple analys	is, reports)	<u> </u>	
E & E/PIT files									
EPA FORM 2070-13 (7-81)									

POTE	ntial hazardous waste site	I. IDENTIFICATION		
EPA	SITE INSPECTION REPORT .	01 STATE 02 SITE NUMBER MO D931715560		
PART 10 -	PAST RESPONSE ACTIVITIES	10 DA3.7.12200		
II. PAST RESPONSE ACTIVITIES				
01A. WATER SUPPLY CLOSED	02 DATE	03 AGENCY		
04_DESCRIPTION				
None known.				
01 B. TEMPORARY WATER SUPPLY PROVIDED	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 C. PERMANENT WATER SUPPLY PROVIDED	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 D. SPILLED MATERIAL REMOVED	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01E. CONTAMINATED SOIL REMOVED	02 DATE	03 AGENCY		
04 DESCRIPTION	- ", "			
None known.				
01 F. WASTE REPACKAGED	02 DATE	03 AGENCY		
04 PESCRIPTION				
None known.				
01 G. WASTE DISPOSED ELSEWHERE	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 x H. ON SITE BURIAL	02 DATE	03 AGENCY		
04 DESCRIPTION				
It was common practice for FMGPs to bury	tar and purifier wastes on site.			
01 I. IN SITU CHEMICAL TREATMENT	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.		•		
01 J. IN SITU BIOLOGICAL TREATMENT	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 K. IN SITU PHYSICAL TREATMENT	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01L. ENCAPSULATION	02 DATE	03 AGENCY		
04 PESCRIPTION				
None known.				
01 M. EMERGENCY WASTE TREATMENT	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01N. CUTOFF WALLS	02 DATE	03 AGENCY		
04 DESCRIPTION -	· ·			
None known.				
01O. EMERGENCY DIKING/SURFACE WATER DI	VERSION 02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 P. CUTOFF TRENCHES/SUMP	02 DATE	03 AGENCY		
04 DESCRIPTION				
None known.				
01 Q. SUBSURFACE CUTOFF WALL	02 DATE	03 AGENCY		
04 DESCRIPTION				
L wone bnown				

	POTENTIAL HAZARDOUS WASTE SITE	I. IDENTIFICATION
EPA	SITE INSPECTION REPORT	01 STATE 02 SITE NUMBER MO D981715980
	PART 10 - PAST RESPONSE ACTIVITIES	557,72500
II. PAST RESPONSE ACTIVITIES (Contin		
01R. BARRIER WALLS CONSTRUCTED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01S. CAPPING/COVERING	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01T. BULK TANKAGE REPAIRED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01U. GROUT CURTAIN CONSTRUCTED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01V. BOTTOM SEALED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		,
01W. GAS CONTROL	02 DATE	03 AGENCY
04 DESCRIPTION None known		
01X. FIRE CONTROL	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01Y. LEACHATE TREATMENT	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01Z. AREA EVACUATED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
01 x 1. ACCESS TO SITE RESTRICTED	(partially) 02 DATE	03 AGENCY
104 DESCRIPTION	ence with barb wire running across the top	surrounds the tank farm area.
012. POPULATION RELOCATED	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
013. OTHER REMEDIAL ACTIVITIES	02 DATE	03 AGENCY
04 DESCRIPTION None known.		
		1
1		
V. SOURCES OF INFORMATION (Cite spe	ecific references, e.g., state files, sample	analysis, reports)
E & E/FIT files		
<i>;</i>		
1		
1		
•		

	POTENTIA	L HAZARDOUS WASTE SITE	I. IDENTIFI	CATION
EPA	SITE	INSPECTION REPORT	01 STATE 02 MO	SITE NUMBER
	PART 11 -	ENFORCEMENT INFORMATION		D381/13480
II. ENFORCEMENT INFORMATION				
01 PAST REGULATORY/ENFORCEMENT ACTION	YES	x_NO		
02 DESCRIPTION OF FEDERAL, STATE, LOC	AL REGULAT	DRY/ENFORCEMENT ACTION		
None known.				
				•
				•
		•		
•	•			
		•		
III. SOURCES OF INFORMATION (Cite s	ecific ref	erences, e.g., state files,	sample analysis, report	.s)
E & E/FIT files			 	

EPA FORM 2070-13 (7-81)

APPENDIX C Site Contacts and Property Owners

Site Contacts

Randel H. Lewis
Terminal Manager
Petroleum, Fuel & Terminal Company
St. Louis, MO 63105
(314) 621-0522

Dick Block Assistant Terminal Manager Terminal Manager Petroleum, Fuel & Terminal Company St. Louis, MO 63105 (314) 621-0522

Bob Welch Terminal Property Manager Terminal Manager Petroleum, Fuel & Terminal Company St. Louis, MO 63105 (314) 621-0522

Greg Reesor
Missouri Coordinator
U.S. Environmental Protection Agency
Region VII
726 Minnesota Avenue
Kansas City, KS
(913) 551-7695

Property Owner Information

Property Owner	Sample #'s
Richard Weidner TRRA 2016 Madison Ave. Granite City, IL 62040	110, 111 201, 202
Marty Anamosa Glacier Park Co. 1011 Western Ave., Ste. 700 Seattle, WA 98104	112
David Bohm City Hall City Counselor's Office, Rm. 314 St. Louis, MO 63103	301 - 303 401 - 403
Doug Mitchem Illinois American Water Co. 300 N. Water Works Dr. Belleville, IL 62223	304, 304D
Randall Lewis PF & T Foot of Mullanphy St. St. Louis, MO 63105	001 - 013 101 - 109 501

APPENDIX D

EPA Data Transmittal

ICF Technology Incorporated NSI Technology Services Corporation The Bionetics Corp.

EPA Region VII Laboratory 25 Funston Rd. Kansas City, KB 66115 (913) 236-3881

To: Larry Marchin & Barry Evans
Data Review Task Monitors

Thru: Harold Brown, Ph.D.

ESAT Deputy Project Officer, EPA

From: Laleh A. Jonooby A. Senior Scientist, ESAT

Thru: Ronald A. Ross

Region VII ESAT Team Leader, NSI-TS

Date: April 19, 1991

Subject: Review of organic data for Laclede Coal Gas

TID #: 07-9103-535 ICF Acct. #: 302-26-535-02

NSI Sales Order: 1073-535 Activity Number: DSX44

ESAT Document Control #:.. ESAT-VII-535-0011

Assignment Number: 764

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses," and the Region VII Organic Data Review Training Manual as guidance. The following comments and attached data sheets are a result of the ESAT review of the above mentioned data from the contract laboratory.

SAS NO.: 6082-G - LABORATORY: RECRA ENV., INC.

SITE: Laclede Coal Gas METHOD NO.: 418.1

REVIEWER: A. Jonooby EPA ACTIVITY NO.: D8X44

MATRIX: Water, Soil and Oil

1			
SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
6082-001	DSX44301	6082-009	DSX44401D
6082-002	DSX44301D	6082-010	DSX44402
6082-003	DSX44302	6082-011	DSX44403
6082-004	DSX44304	6082-012	DSX44501
6082-005	DSX44304D		
60 82- 00 6	DSX44207F		
6082-007	DSX44303		
6082-008	DSX44401		

And QC samples DSX44304 (MS), 303 (DUP), 401 (DUP), 401D (MS), 501 (MS), 501 (DUP), 910M, 911M and 912M.

General

This data review assignment covers <u>SEVEN WATER</u> samples (DSX44301 - 304 and 207F), <u>FOUR SOIL</u> samples (DSX44401 - 403) and <u>ONE OIL</u> sample (DSX44501) analyzed for total petroleum hydrocarbons (TPHs). Field blank and field duplicate were included in this assignment. TPHs traffic reports were present.

1. Holding Times and Preservation

There are no established holding times for TPHs. Sample preservation is to keep cold at 4 degrees C.

2. Method Blanks / Field Blanks

All of the method blanks and field blank were reported below method detection limits. For the water samples, the field blank and the method blank were reported as <0.5 mg/L, for soil samples, the method blank was reported to be <25 mg/L and the method blank analyzed with the oil sample was reported to be <1000 mg/L.

3. Initial Calibrations

A six point calibration curve using reference oil #1 as the source at an absorbance of 2930 cm-1 was established for this activity. A correlation coefficient of 0.9999 was achieved with this curve.

4. Matrix Spike

A matrix spike was run with each set of matrices. Percent recovery of the MS for the water samples was 84.8%, for soils, the MS recovery was 96.8% and the MS recovery for the oil sample was 105.5%.

5. Lab Duplicate

A lab duplicate was run with each set of matrices. The relative percent difference between the water sample and its duplicate was not calculated because they were reported to be less than the method detection limits. A 15% relative percent difference was calculated for the soil samples and 1.2% relative percent difference was reported for the oil sample.

Discussion

Sample DSX44301 was broken accidently by RECRA Environmental lab before any TPH analysis could be done on it. Therefore, there is no data available for this sample in this package. However, RECRA Environmental, Inc. did contact SMO with regards to this matter as mentioned in their memo and since there was a duplicate for that sample (DSX44301D) no action was taken by the EPA to resample.

All of the water samples were extracted using 1L from each sample. The TPHs were reported less than the method detection limits (<0.5 mg/L).

A sample weight of 20g was used for the soil samples and only a 0.5g was used for the oil sample with a dilution factor of 50.

Summary

From the evaluation of the QC mentioned above, this data meets the requirements for overall accuracy, precision and completeness.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

ICF Technology Incorporated

NBI Technology Services Corp.

The Bionetics Corp.

ESAT Region VII NSI Technology Services 25 Funston Road Kansas City, KS 66115 (913) 236-3881

MEMORANDUM

TO: Larry Marchin, Chemist, CLQA/LABO/ENSV

THRU: Harold Brown, Ph.D, DPO/LABO/ENSV

FROM: Peggy Cox, QA/QC Chemist/ESAT/MANTECH Q OFFICE

THRU: Ronald Ross, Manager/ESAT/MANTECH

DATE: April 27, 1991

SUBJECT: Review of data for LACLEDE COAL GAS

TID#: 07-9103-535

ASSIGNMENT#: 760

ICF ACCT#: 302-26-535-02

MANTECH S.O.#: 1073-535

ESAT DOCUMENT#: ESAT-VII-535-0008

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," July 1, 1988 revision.

The following comments and attached data sheets are a result of ManTech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE/SAS NO.: 15978

CONTRACT NO.: 68-D0-0108

SITE: LACLEDE COAL GAS

LABORATORY: SKINER

METHOD NO.: CS03901

EPA ACTIVITY: DSX44

REVIEWER: P. COX MATRIX: WATER

SMO SAMPLE NO. EPA SAMPLE NO.

MGH515 DSX44201 **MGH517** DSX44202 MGH519 DSX44301 MGH521 DSX44301D MGH523 DSX44302 **MGH525** DSX44304 ' MGH527 DSX44304D MGH529 DSX44900P

GENERAL

Case 15978 contained 8 environmental and 6 QC water samples analyzed for total metals and cyanide at the low level concentration. No field blank, two field duplicates, and one performance evaluation sample were included in this data review assignment. Data review was performed at level 2.

1. TECHNICAL HOLDING TIMES and PRESERVATION

A. Technical holding times and required preservation were within quality control limit requirements.

2. INITIAL and CONTINUING CALIBRATION

A. Initial and continuing calibrations were within quality control limit requirements.

3. BLANKS

- A. No analytes were detected above the contract required detection limit (CRDL) in any blank.
- B. Levels of analytes greater than the instrument detection limit (IDL) were detected for aluminum (Al), barium (Ba), calcium (Ca), cadmium (Cd), iron (Fe), nickel (Ni), silver (Ag), and zinc (Zn). Analytes greater than the instrument detection limit (IDL) but less than 5 times the highest level detected in the blank were qualified with a "U" code. Cadmium (Cd) in sample DSX44900P and zinc (Zn) in samples DSX44301, DSX44301D, and DSX44302 were qualified with a "U" code according to the blank rules.

4. ICP INTERFERENCE CHECK

A. All analytes contained in the ICP interference check sample were within quality control limit requirements. Antimony (Sb), potassium (K), and sodium (Na) were found but not elements present in the AB ICP interference check solution. Since levels detected were below the intrument detection limit (IDL) or levels found in the samples were significantly higher, no data were qualified by the ICP interference check sample.

5. LABORATORY CONTROL SAMPLE

A. All laboratory control samples analyzed were within quality control limit requirements.

6. DUPLICATES

A. All analytes were within quality control limit requirements except aluminum (Al) and cyanide (CN⁻). Aluminum (Al) in samples DSX44201, DSX44202, DSX44301, DSX44301D, DSX44302, DSX44304, DSX44304D, DSX44202S, and DSX44901C and cyanide (CN⁻) in samples DSX44201, DSX44202, DSX44304D, DSX44900P, and DSX44202S were qualified with a "J" code according to the duplicate rules.

7. MATRIX SPIKE

A. All analytes were within quality control limit requirements for percent recovery except selenium (Se) and thallium (Tl). Selenium had a 0.0% recovery due to the original sample being diluted. No samples were qualified for selenium (Se) but thallium (Tl) in sample DSX44901C was qualified with a "J" code according to the spike rules.

8. GRAPHITE FURNACE ATOMIC ABSORPTION (GFAA) SPECTROSCOPY

- A. All quality control limit requirement criteria for graphite furnace atomic absorption (GFAA) spectroscopy were met. Selenium (Se) and thallium (Tl) in several samples had post digestion spike recoveries outside quality control limits, but since the results were non-detect, no data were qualified.
- B. The method of standard additions was performed for arsenic (As) and selenium (Se) in sample DSX44900P and lead (Pb) in sample DSX44202. The correlation coefficient for lead (Pb) in sample DSX44202 and selenium (Se) in sample DSX44900P was outside quality control limit requirements (less than 0.995) and were qualified with a "J" code according to the standard addition rules.

9. PERFORMANCE EVALUATION AUDIT SAMPLE

A. Performance evaluation audit sample DSX44900P was submitted to the laboratory for analysis with all analytes containined in the audit being identified.

10. ICP SERIAL DILUTION

A. All analytes were within quality control limit requirements except copper (Cu). Copper (Cu) in samples DSX44201, DSX44202, DSX44202S, DSX44201L, and DSX44901C were qualified with a "J" code according to the ICP serial dilution rules.

11. SUMMARY

- A. Cadmium (Cd) in sample DSX44900P and zinc (Zn) in samples DSX44301, DSX44301D, and DSX44302 were qualified with a "U" code due to blank contamination.
- B. Aluminum (Al) in samples DSX44201, DSX44202, DSX44301, DSX44301D, DSX44302, DSX44304, DSX44304D, DSX44202S, and DSX44901C and cyanide (CN) in samples DSX44201, DSX44202, DSX44304D, DSX44900P, and DSX44202S were qualified with a "J" code due to poor duplicate precision.
- C. Thallium (T1) in sample DSX44901C was qualified with a "J" code due to poor spike recovery.
- D. Lead (Pb) in sample DSX44202 and selenium (Se) in sample DSX44900P were qualified with a "J" code according to the standard addition rules.
- E. Copper (Cu) in samples DSX44201, DSX44202, DSX44202S, DSX44201L, and DSX44901C were qualified with a "J" code according to the ICP serial dilution rules.
- F. Several compounds were "U" coded due to the compound concentrations being greater than the instrument detection limit (IDL) but less than the contract required detection limit (CRDL).
- G. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Inorganic Analysis dated March 1990, with the exceptions noted above.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

ICF Technology Incorporated

NSI Technology Services Corp.

The Bionetics Corp.

ESAT Region VII NSI Technology Services 25 Funston Road Kansas City, KS 66115 (913) 236-3881

MEMORANDUM

TO:

Larry Marchin, Chemist, CLQA/LABO/ENSV

THRU:

Harold Brown, Ph.D, DPO/LABO/ENSV

FROM:

Peggy Cox, QA/QC Chemist/ESAT/MANTECH & Coffekt

THRU:

Ronald Ross, Manager/ESAT/MANTECH

DATE:

SUBJECT: Review of data for LACLEDE COAL GAS

TID#:

07-9103-535

ASSIGNMENT#:

761

ICF ACCT#:

302-26-535-02

MANTECH S.O.#: 1073-535

ESAT DOCUMENT#: ESAT-VII-535-0010

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic

Analyses," July 1, 1988 revision.

The following comments and attached data sheets are a result of ManTech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE/SAS NO.: 15978

LABORATORY:

SKINER

CONTRACT NO.: 68-D0-0108

METHOD NO.: CS0390I

SITE: LACLEDE COAL GAS

EPA ACTIVITY: DSX44

REVIEWER: P. COX

MATRIX: WATER

SMO SAMPLE NO.

EPA SAMPLE NO.

MGH516	
MGH518	
MGH520	
MGH522	
MGH524	
MGH526	

MGH528

MGH530

DSX44201 DSX44202 DSX44301 DSX44301D DSX44302 DSX44304

DSX44304D DSX44900P

GENERAL

Case 15978 contained 8 environmental and 6 QC water samples analyzed for dissolved metals at the low level concentration. No field blank, two field duplicates, and one performance evaluation sample were included in this data review assignment. Data review was performed at level 2.

1. TECHNICAL HOLDING TIMES and PRESERVATION

A. Technical holding times and required preservation were within quality control limit requirements.

2. INITIAL and CONTINUING CALIBRATION

A. Initial and continuing calibrations were within quality control limit requirements.

3. BLANKS

- A. No analytes were detected above the contract required detection limit (CRDL) in any blank.
- B. Levels of analytes greater than the instrument detection limit (IDL) were detected for barium (Ba), calcium (Ca), copper (Cu), and zinc (Zn). Analytes greater than the instrument detection limit (IDL) but less than 5 times the highest level detected in the blank were qualified with a "U" code. Since all associated data were non-detect, no data were qualified by the blank rules.

4. ICP INTERFERENCE CHECK

A. All analytes contained in the ICP interference check sample were within quality control limit requirements. Antimony (Sb), potassium (K), and sodium (Na) were found but not elements present in the AB ICP interference check solution. Since levels detected were below the intrument detection limit (IDL) or levels found in the samples were significantly higher, no data were qualified by the ICP interference check sample.

5. LABORATORY CONTROL SAMPLE

A. All laboratory control samples analyzed were within quality control limit requirements.

6. DUPLICATES

A. All analytes were within quality control limit requirements except lead (Pb). Lead (Pb) in samples DSX44201, DSX44301D, DSX44900P, DSX44201S, and DSX44903C were qualified with a "J" code according to the duplicate rules.

7. MATRIX SPIKE

A. All analytes were within quality control limit requirements for percent recovery except selenium (Se) and thallium (T1). Selenium (Se) in samples DSX44900P and DSX44903C and thallium (T1) in sample DSX44903C were qualified with a "J" code according to the spike rules.

8. GRAPHITE FURNACE ATOMIC ABSORPTION (GFAA) SPECTROSCOPY

- A. All quality control limit requirement criteria for graphite furnace atomic absorption (GFAA) spectroscopy were met. Selenium (Se) and thallium (T1) in several samples had post digestion spike recoveries outside quality control limits, but since the results were non-detect, no data were qualified.
- B. The method of standard additions was performed for arsenic (As) in sample DSX44202. The correlation coefficient for arsenic (As) in sample DSX44202 was outside quality control limit requirements (less than 0.995) and was qualified with a "J" code according to the standard addition rules.

9. PERFORMANCE EVALUATION AUDIT SAMPLE

A. Performance evaluation audit sample DSX44900P was submitted to the laboratory for analysis with all analytes containined in the audit being identified.

10. ICP SERIAL DILUTION

A. All analytes were within quality control limit requirements except iron (Fe). Iron (Fe) in samples DSX44201, DSX44202, DSX442015, DSX44201L, and DSX44903C were qualified with a "J" code according to the ICP serial dilution rules.

11. SUMMARY

A. Lead (Pb) in samples DSX44201, DSX44301D, DSX44900P, DSX44201S, and DSX44903C were qualified with a "J" code due to poor duplicate precision.

- B. Selenium (Se) in samples DSX44900P and DSX44903C and thallium (T1) in sample DSX44903C were qualified with a "J" code due to poor spike recovery.
- C. Arsenic (As) in sample DSX44202 was qualified with a "J" code according to the standard addition rules.
- D. Iron (Fe) in samples DSX44201, DSX44202, DSX44201S, DSX44201L, and DSX44903C were qualified with a "J" code according to the ICP serial dilution rules.
- E. Several compounds were "U" coded due to the compound concentrations being greater than the instrument detection limit (IDL) but less than the contract required detection limit (CRDL).
- F. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Inorganic Analysis dated March 1990, with the exceptions noted above.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- ZONE II

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Larry Marchin & Barry Evans To:

Data Review Task Monitors

Thru: Harold Brown, Ph.D.

ESAT Deputy Project Officer, EPA

From: Janet Muse

Senior Sclentist, ESAT

Thru: Ronald A. Ross

Region VII ESAT Team Leader. Mantech

Date: April 26, 1991

Subject: Review of organic data for Laclede Coal Gas

ASGN #: 777

TID #: 07-9103-535 ICF Acct. #: 302-26-535-02

ManTech Sales Order: 1073-535

ESAT Doc. Control #:.... ESAT-VII-535-0024

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses," and the Region VII Organic Data Review Training Manual as guidance. The following comments and attached data sheets are a result of the ESAT review of the above mentioned data from the contract laboratory.

CASE NO.: 15978

SITE: Laclede Coal Gas

REVIEWER: J. MUSE

MATRIX: Oil

LABORATORY:

EIRA

METHOD NO.: CS03900 EPA ACTIVITY NO.: DSX44

Sample No. GJ894

EPA Sample No. DSX44501

And QC samples DSX44919M, 919G, and 919H.

General

This data review assignment covers **ONE OIL** sample analyzed for High Concentration volatile organics and extractables SOW 9/88. No field blanks or field duplicates, and 4 QC samples were included in this assignment. The organic traffic report was present.

Holding Times and Preservation

There are no holding time requirements for volatile organics in oil samples.

K HI. Cone

Preservation of samples was refrigeration at 4 degrees C and protection from light.

2. GC/MS Performance

All CLP criteria for GC/MS performance were met for a 50 nanogram injection of 4-bromofluorobenzene.

All CLP criteria for GC/MS performance were met for a 50 nanægram injection of DFTPP.

3. Method Blanks

Positive hits for target compounds in the blanks were only reported positive in the sample after the blank rule was applied. That is, positive hits for methylene chloride in the sample were reported and coded with a "U" unless they were greater than 10-times the level in the daily blank. Methylene chloride was reported in the VOC method blank. No target compounds for the extractables were found in the method blank.

4. Initial Calibration

VOCS: All response factors met applicable criteria. All compounds met the 35% criteria for RSD.

Extractables: A three point initial calibration was done before analysis of the samples. All response factors met applicable criteria. All compounds met the 35% criteria for RSD.

5. Continuing Calibration

VOCS: The sample and QC samples were analyzed the same day of the initial calibration, so no continuing calibration std. needed to be run.

Extractables: All compounds were within the 25% difference criteria for response factors on the continuing calibration of 3/29/91, except for a few target compounds. Since no positive hits were found for these outliers in the sample, no data was qualified due to the continuing calibration.

6. Surrogate Recoveries

VOCS: Surrogate recoveries were within control limits in all cases.

Extractables: Surrogates were diluted out of the sample due to large quantities of target compounds. No data were qualified due to this.

7. Internal Standard Areas

VOCS: Met applicable criteria.

Extractables: Perylene-d12 and chrysene-d12 were less then the lower limit criteria in the method blank. No data was qualified due to this.

8. Control Matrix Spike

A control matrix spike was done for both parameters, VOCS and extractables. Recoveries were within control limits for both control matrix spikes.

Summary

This data package mets acceptability in terms of the requirements for overall accuracy, precision and completeness.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

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ManTech Environmental Technology, Inc.

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TO:

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THRU:

Harold Brown, Ph.D., ESAT Contract Manager/ENSV

FROM:

Rebecca K. Estep, ESAT Data Reviewer/ManTech

THRU:

Albert Iannacone, ESAT QA/QC Coordinator/ManTech

DATE:

April 29, 1991

SUBJECT: Review of organic data for Laclede Coal Gas.

TID#:

07-9103-535

ASSIGNMENT#:

769

ICF ACCT#:

302-26-535-02

ManTech S.O.#:

1073-535

ESAT Document#:

ESAT-VII-535-0014

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses," February 1, 1988 revision.

The following comments and attached data sheets are a result of ManTech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE NO.:

15978

LABORATORY:

RECMD

CONTRACT NO.: SITE:

68-D0-0158

METHOD NO: EPA ACTIVITY: CS03900 DSX44

REVIEWER:

Laclede Coal Gas Rebecca K. Estep

MATRIX:

Soil

Samples for Semivolatile Analysis

SMO SAMPLE NO.	EPA SAMPLE NO.	SMO SAMPLE NO.	EPA SAMPLE NO.
G J8 79	DSX44108	GX983	DSX44002
GJ880	DSX44109	GK984	DSX44003
*GJ890	*DSX44401	GK985	DSX44004
*GJ891	*DSX44401D	GK986	DSX44005
*GJ892	*DSX44402	GK987	DSX44110
*GJ893	*DSX44403	GK988	DSX44111
GK982	DSX44001	GK989	DSX44112

^{*} Also analyzed for volatiles.

GENERAL

Case 15978 contained low level soil samples for volatile and base/neutral and acid analyses. 14 samples for semivolatile analysis, 4 which were also analyzed for volatiles (18 actual;13 QC). This package includes no field blanks, one field duplicate, and no performance evaluation sample. Data review was performed at level 2.

1. TECHNICAL HOLDING TIMES and PRESERVATION

A. No technical holding times are specified for soil samples. All samples were within holding time specifications from extraction to analysis for base/neutral and acids.

2. GC/MS Tuning

- A. All volatile GC/MS tunings and mass calibrations were within quality control limit requirements for bromofluorobenzene (BFB).
- B. All base/neutral and acid GC/MS tunings and mass calibratons were within quality control limit requirements for decafluorotriphenylphosphine (DFTPP).

3. INITIAL and CONTINUING CALIBRATION

- A. All average relative response factors and percent relative standard deviations were within quality control limit requirements for the volatile initial calibration.
- B. Bromoform and 4-methyl-2-pentanone were outside quality control limit requirements for percent difference (%D) (greater than 25%) on the volatile continuing calibration dated 3/14/91. Since these compounds were non-detect in all samples no data were qualified by the calibration rules.
- C. All average relative response factors and percent relative standard deviations were within quality control limit requirements for the base/neutral and acid initial calibrations except for the one dated 4/8/91 where 4-chlorophenyl-phenyl ether, fluorene, and bis(2-ethylhexyl)phthalate were outside the quality control limit requirements for percent relative standard deviation (%RSD) (greater than 30%). This resulted in fluorene being "J" coded in sample DSX44109, and bis(2-ethylhexyl)phthalate "J" coded in samples DSX44001, DSX44002, DSX44003, DSX44004, and DSX44005.

D. Several compounds (anywhere from 2 to 7) were outside quality control limit requirements for percent difference (%D) (greater than 25%) on the base/neutral and acid continuing calibrations. Pyrene was "J" coded in sample DSX44402 from the continuing calibration dated 4/3/91. Bis(2-ethylhexyl)phthalate would have been "J" coded in samples DSX44001, DSX44002, DSX44003, DSX44004, and DSX44005 from the continuing calibration dated 4/10/91, however this compound was previously "J" coded in these same 5 samples due to the initial calibration rules.

4. INTERNAL STANDARD RESPONSE

- A. All internal standard response areas were within a factor of 2 when comparing samples and their associated continuing calibration response areas for volatile analysis.
- B. The internal standard response areas were out by more than a factor of 2 between the samples and their associated continuing calibration response areas for base/neutral and acid analysis in sample DSX44108 and the reanalysis of DSX44108 for phenanthrene-d10 and chrysene-d12. DSX44108 was also out for perylene-d12. This resulted in carbazole being "J" coded for sample DSX44108.

5. BLANKS

- A. One method blank was analyzed for the volatile fraction with acetone, methylene chloride, and 2-butanone detected. This resulted in methylene chloride and acetone being "U" coded in samples DSX44401D, DSX44402, DSX44401S, and DSX44401W, and only acetone in sample DSX44403 according to the blank rules. Methylene chloride and 2-butanone, in the method blank, were detected above the instrument detection limit (IDL) but less than the contract required quantitaion limit (CRQL), thus their results were raised to the CRQL and "U" coded.
- B. Four method blanks were analyzed for the base/neutral and acid fraction with bis(2-ethylhexyl)phthalate detected in all 4 samples. No qualifications were made because the concentrations in all corresponding samples were greater than ten times the blank concentrations. In addition, butylbenzylphthalate was detected in sample DSX44924M and DSX44925M, and diethylphthalate and di-n-butylphthalate were detected in sample DSX44926M. Since these compounds were non-detect no data were qualified due to the blank rules.

6. SURROGATE RECOVERY

A. All surrogates were within quality control limit requirements for percent recovery for the volatile fraction except 1,2-dichloroethane-d4 in sample DSX44401W. No data were qualified due to the surrogate recovery rules.

B. All surrogates were within quality control limit requirements for percent recovery for the base/neutral and acid fraction except 2-fluorobiphenyl and terphenyl-d14 in sample DSX44108DL. Phenanthrene, fluoranthene, pyrene, benzo(a)anthracene, chrysene, benzo(b)fluoranthene, benzo(a)pyrene, indeno(1,2,3-cd) pyrene, and benzo(g,h,i)perylene were "J" coded in sample DSX44108 due to the fact that the concentrations for these base/neutral compounds were reported from the dilution and the dilution was qualified due to the surrogate rules. Terphenyl-d14 was also out for samples DSX44108 and DSX44108RE, however no data were qualified.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

- A. A matrix spike/matrix spike duplicate was analyzed for volatiles with percent recovery and relative percent difference within quality control limit requirements.
- B. A matrix spike/matrix spike duplicate was analyzed for the base/neutral and acid fraction with the percent recovery for pyrene low and the relative percent difference for acenaphthene and pyrene outside the advisory limits. No data were qualified due to the matrix spike/matrix spike duplicate rules.

8. PERFORMANCE EVALUATION SAMPLE

A. No performance evaluation sample was submitted to the laboratory for analysis associated with this case and SDG number.

9. COMPOUND IDENTIFICATION and QUANTITATION

- A. All target compound identifications were supported by good agreement between sample and standard mass spectra submitted for each positive sample response.
- B. Several compounds in all samples for both volatile and base/neutral and acid analyses were detected above the instrument detection limit (IDL) but below the contract required quantitation limit (CRQL). These results were raised to the CRQL and "U" coded.

10. SUMMARY

- A. Samples DSX44401D, DSX44402, DSX44403, DSX44401S, and DSX44401W for volatile analysis detected methylene chloride and/or acetone which were "U" coded due to the blank rule.
- B. Anthracene was "J" coded in sample DSX44108 due to the high concentration encountered in both the original analysis and the dilution which were detected above the calibration range.
- C. Fluorene was "J" coded in sample DSX44109, and bis(2-ethylhexyl)phthalate was "J" coded in samples DSX44001, DSX44002, DSX44003, DSX44004, and DSX44005 due to the initial calibration rules.

- D. Pyrene was "J" coded in sample DSX44402 due to the continuing calibration rules.
- E. Carbazole was "J" coded in sample DSX44108 due to the internal standard rules.
- F. Several compounds in all samples for volatile and base/neutral and acid analyses were detected above the IDL but less than the CRQL, thus these compounds were raised to the CRQL and "U" coded.
- G. Phenanthrene, fluoranthene, pyrene, benzo(a)anthracene, chrysene, benzo(b)fluoranthene, benzo(a)pyrene, indeno(1,2,3-cd) pyrene, and benzo(g,h,i)perylene were "J" coded in sample DSX44108 due to the surrogate rules.
- H. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Organic Analysis dated March 1990, with the exceptions noted above.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM -- Zone II.

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TO:

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Data Review Task Monitors

THRU:

Harold Brown, Ph.D.

ESAT Deputy Project Officer, EPA

FROM:

Paula Woodland Daw (

ESAT Scientist Y

THRU:

Ronald A. Ross

ESAT Team Manager

DATE:

April 29, 1991

SUBJECT: Review of VOA/BNA data for Laclede Coal Gas

TID# 07-9103-535 ASSIGNMENT# 768 ICF ACCT# 26-535-02 ManTech S.O.# 1073-535

ESAT Document #: ESAT-VII-535-0013

These data were reviewed primarily according to the Special Analytical Services Request and the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses," February 1988 revision with changes given in the Region VII Organic Data Review Training Manual and EPA memorandums and the 3/90 SOW.

The following comments and attached data sheets are a result of the ESAT review, according to EPA policies, of the following data from the contract laboratory.

CASE NO.: 15978
SITE: Laclede Coal Gas
REVIEWER: Paula Woodland

LABORATORY: RECMD
METHOD NO.: CS03900
EPA ACTIVITY NO.: DSX44

MATRIX: SOIL

<u>VOA</u>

SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
GJ859	DSX44001	GJ865	DSX44007
GJ860	DSX44002	GJ866	DSX44008
GJ861	DSX44003	GJ867	DSX44009
GJ862	DSX44004	GJ868	DSX44010
GJ863	DSX44005	GJ869	DSX44011
GJ864	DSX44006	GJ870	DSX44012
		GTR71	DSX44013

SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
GJ864	DSX44006	GJ872	DSX44101
GJ865	DSX44007	G J87 3	DSX44102
GJ866	DSX44008	GJ874	DSX44103
G J86 7	DSX44009	G J87 5	DSX44104
G J8 69	DSX44011	G J8 76	DSX44105
GJ870	DSX44012	G J87 7	DSX44106
GJ871	DSX44013	G J8 78	DSX44107

GENERAL

This data review assignment covers 12 <u>SOIL</u> samples analyzed for <u>VOA</u> and 14 <u>SOIL</u> samples analyzed for <u>BNA</u>. There were no field blanks, duplicates, or performance evaluation samples included with this assignment. DSX44010 (GJ868) was originally sent along with this SDG number, however SMO penoved the sample to another SDG number because the matrix was defined incorrectly.

1. Holding Times and Preservation

VOA: Technical holding times have not been established for soil samples.

BNA: Collection to extraction holding times have not been established for soil samples, however the extraction to analysis holding times were within the established criteria.

2. GC/MS Tuning

All relative ion abundances were within the established control limits of the 3/90 SOW for both VOA and BNA analysis.

3. Initial and Continuing Calibration

VOA: The %RSD was out of control for acetone in the initial calibration curve ran on 1/8/91. All positive results for this compound in samples DSX44005 and DSX44913M were coded J. All criteria were within the established control limits for the other initial calibration curve. The %D was out of control for several compounds in the continuing calibration check, however no other samples were coded.

BNA: All criteria were within established control limits for the initial calibration curve. The %D was out of control for pyrene and several other compounds on 3-22-91, 3-26-91, and 3-28-91. All positive results for pyrene in samples DSX44007, -009, -011, -1015, and -101W were coded J.

4. Internal Standard Response

VOA: The internal standard area was out of control for chlorobenzene-d⁵ in sample DSX44012 and the re-analysis of DSX44012. The positive result for the corresponding compound, styrene, in DSX44012 was coded J. The original analytical result was reported because the internal standard response was slightly better than the re-analysis. All other internal standard criteria were within control limits.

433

BNA: The internal standard area was out of control for naphthalene-d⁸ in sample DSX44008 and the re-analysis, but only non-detect compounds were reported from these injections so no coding was performed. The internal standard response was within the established criteria for the dilution of DSX44008 so the diluted results for naphthalene and 2-methylnaphthalene were not coded. All other internal standard criteria were within control limits.

5. Blanks

VOA: Methylene chloride and acetone were detected in the blanks. Corresponding sample results were qualified according to the blank rule using ten times the highest blank value. Acetone was qualified in samples DSX44001, -002, -003, -004, -012, and -001S. Methylene chloride was qualified in samples DSX44001, -002, -001S, and -001W.

BNA: Bis(2-ethylhexyl)phthalate was detected in the blanks, however no qualifications were performed.

6. Surrogate Recovery

The surrogate compound recoveries were within established control limits.

7. Matrix Spike/Matrix Spike Duplicate Recovery

VOA: Matrix spike/matrix spike duplicate criteria were within established control limits.

BNA: The % recovery of the medium level soil matrix spike and matrix spike duplicate was out of control for 6 and 2 of the 11 compounds, respectively. The %RPD was out of control for 3 of the 11 compounds. No coding was performed. A low level matrix spike and matrix spike duplicate from a different SDG# (GJ879) was included in this SDG# GJ859 for the one low level sample (DSX44106). This was not evaluated because it was from another case.

8. Compound Identification and Quantitation

Due to the requested review level, results listed on the summary forms were used for the review. These results were not checked against the raw data for accuracy, and calculations were not verified.

VOA: Xylene was coded J in sample DSX44008 because it was detected over the calibration range. A re-analysis was not performed by the lab.

BNA: For DSX44008, nine compounds were over the calibration range so a dilution was made. In this dilution, seven of the nine compounds were below the CRQL. These seven compounds were reported using the original undiluted results and coded J for exceeding the calibration range. The seven compounds that were coded J in DSX44008 were fluorene, phenanthrene, anthracene, fluoranthene, pyrene, benzo(a)anthracene, and chrysene.

9. Summary

VOA: Acetone was coded J for sample DSX44012 because the internal standard response was too low. Also, acetone was coded J for samples DSX44005 and -913M because of an initial calibration RSD outlier. Xylene was coded J for sample DSX44008 because it was over the calibration range. Several samples were qualified by the blank rule for methylene chloride and acetone.

BNA: Pyrene was coded J for samples DSX44007, -009, -011, -101S, and -101W because the %D was outside of control limits for the corresponding continuing calibration standard. An improper dilution was made for sample DSX44008 which resulted in the coding of several compounds (see Compound Identification and Quantitation section above). A low level soil matrix spike and matrix spike duplicate was not included with this SDG number.

The EPADA data sheets are attached. This data review is complete. If you have any questions, please contact me.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II.

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TO:

Larry Marchin/Barry Evans Data Review Task Monitor

THRU:

Harold Brown, Ph.D.

ESAT Deputy Project Officer, EPA

FROM:

Kevin Ludwikoski

ESAT Data Reviewer

THRU:

Ronald A. Ross ESAT Team Manager

DATE:

April 29, 1991

Review of inorganic data for Laclede Coal Gas Co. SUBJECT:

> TID# 07-9103-535 ASSIGNMENT# 762

ICF ACCT# 302-26-535-02

ManTech S.O.# 1073-535 ESAT Document No. ESAT-VII-535-0025

These data were reviewed primarily according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," July 1988 revision with changes given in the Region VII Inorganic Data Review Training Manual and EPA memorandums.

The following comments and attached data sheets are a result of the ESAT review, according to EPA policies, of the following data from the contract laboratory.

CASE NO.: 15978

SITE: Laclede Coal Gas Co.

REVIEWER: Kevin Ludwikoski

LABORATORY: METHOD NO.:

SKINER

CS0390I EPA ACTIVITY NO.: DBX44

MATRIX: WATER/SOIL

TOTAL METALS and CN /COTT \

	(201	. Lu }	
SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
MGG558	DSX44006	MGG568	DSX44103
MGG559	DSX44007	MGG569	DSX44104
MGG560	DSX44008	MGG570	DSX44105
MGG561	DSX44009	MGG571	DSX44106
MGG562	DSX44010	MGG572	DSX44107
MGG563	DSX44011	MGG573	DSX44108
MGG564	DSX44012	MGG574	DSX44109
MGG565	DSX44013	MGG590	DSX44401
MGG566	DSX44101	MGG591	DSX44401D
MGG567	DSX44102	MGG592	DSX44402

DISSOLVED METALS (WATER)

SMO Sample No.	EPA Sample No.	
MGG576	DSX44203	
MGG578	DSX44204	
MGG583	DSX44206D	
MGG585	DSX44207F	
MGG587	DSX44208	
MGG589	DSX44303	

GENERAL

This data review assignment covers <u>SIX WATER</u> samples analyzed for <u>DISSOLVED METALS</u> and <u>TWENTY SOIL</u> samples analyzed for <u>TOTAL</u> <u>METALS and CYANIDE</u> for case number <u>15978</u>. There was one field blank and one field duplicate for Dissolved Metals and one field duplicate for Total Metals and Cyanide included with this assignment.

1. Technical Holding Times / Preservation

- A. Technical holding times were observed for all analytes.
- B. Technical holding times are not specified for soil samples.

2. Initial and Continuing Calibration

- A. All percent recoveries were within control limits for water samples.
- B. All percent recoveries were within control limits for soil samples.

3. Blanks

A. Several analytes were detected in the blanks. Corresponding sample results were qualified according to the blank rule using <u>five times</u> the highest blank value. Sample results requiring modification are reported as ron-detect on the attached data sheets.

DISSOLVED METALS (WATER)

<u>Analyte</u>	Blank (uq/L)	Qualified Samples
Al	186	None qualified
Ca	1360	None qualified
Cn	24	None qualified
Fe	180	D8Y44208

5 x Highest

. —	400	DDA442U0
Pb	11	DSX44203, -204, -206D, -203L
Mg	372	None qualified
Na	1055	None qualified
V	14	None qualified
Zn	49	None qualified
Mn	9	None qualified

3. Blanks (cont.)

B. TOTAL METALS and CN (SOILS)

Analyte	5 x Highest <u>Blank (mg/kg</u>)	Oualified Samples
Al	30	None qualified
Sb	20	None qualified
Ba	2.4	None qualified
Ca	60	None qualified
Co	3.3	None qualified
Cu	2.8	None qualified
Fe	26	None qualified
Mg	66	None qualified
Mn	6.6	None qualified
, Na	78	None qualified
Zn	3.3	None qualified

4. ICP Interference Check

- A. Recoveries of solution AB analytes were within control limits for Dissolved Metals. Sb, K, and Na were found but not present in the ICS solution. All absolute values for these analytes were below the CRDL and IDL; therefore, no data were qualified by the ICP interference check sample.
- B. Recoveries of solution AB analytes were within control limits for Total Metals and Cyanide. Sb, K, and Na were found but not present in the ICS solution. All absolute values for these analytes were below the CRDL and IDL; therefore, no data were qualified by the ICP interference check sample.

5. Laboratory Control Standard (LCS)

- A. LCS results were within established control limits for Dissolved Metals.
- B. LCS result: were within established control limits for Total Metals and CN

6. <u>Duplicates</u>

- A. The RPD for all analytes were within control limits for Dissolved Metals.
- B. The RPD for Pb exceeded control limits for Total Metals. The sample results that were coded J are listed below.

TOTAL METALS and CN (SOIL)

<u>Analyte</u>	Samples Oualified	<u>Code</u>
Pb	DSX44006-013, DSX44101-109, -401, -401D -4015, -402, -922C	J

7. Matrix Spike Sample

A. Se and Tl were out of range for matrix spike recovery for Dissolved Metals. The samples that had data qualified are listed below.

DISSOLVED METALS

Analyte % Rec	(WATER) Sample No.	Code
Se 45.6	DSX44303, -921C	J
Tl 34.1	DSX44921C	J

B. Sb was out of range for matrix spike recovery for Total Metals. The samples that had data qualified are listed below.

TOTAL METALS and CN (SOIL)

<u>Analyte</u>	<pre>% Rec</pre>	Sample No.	<u>Code</u>
Sb	37.5	DSX44922C	J

8. ICP Serial Dilution

A. The ICP serial dilution results for Zn were outside control limits for Dissolved Metals. The samples that had results qualified are listed below.

DISSOLVED METALS (WATER)

<u>Analyte</u>	Samples Oualified	<u>Code</u>
Zn	DSX44203, -204, -206D, -204S, -203L, -921C	J

B. The ICP serial dilution results for Zn were outside control limits for Total Metals. The samples that had results qualified are listed below.

TOTAL METALS and CN (SOIL)

<u>Analyte</u>	Samples Oualified	Code
Zn	DSX44006, -007, -008, -009, -010, -011, -012, -013, -101, -102, -103, -104, -105, -106, -107, -108, -109, -401, -401D, -401L, -401S, -402, -922C	J

9. Purnace Atomic Absorption

- A. The analytical scheme was followed for Furnace AA analysis for Dissolved Metals. MSA correlation coefficients were acceptable for As in sample DSX44206D (cc=.9987) and for Se in sample DSX44303 (cc=.9964).
- B. The analytical scheme was followed for Furnace AA analysis for Total Metals. MSA correlation coefficients were not acceptable for Se in samples DSX44103, -401L, -401D, and -402 (best cc=.9869, .9878, .9565, and .9800 respectively). Only sample DSX44402 had Se found above the CRDL, and was "J" coded accordingly.

10. Summary

- A. Nine analytes were qualified by the blank rule, one by the the spike rule, and one analyte was qualified by serial dilution outliers.
- B. Eleven analytes were qualified by the blank rule, one by the spike rule, one by the duplicate rule, one by the serial dilution rule, and one analyte in one sample was "J" coded due to the rules of standard additions.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II.

ICF Technology Incorporated

NBI Technology Services Corp.

The Bionetics Corp.

ESAT Region VII

NSI Technology Services

25 Funston Road

Kansas City, KS 66115

(913) 236-3881

TO:

Larry Marchin/Barry Evans

Data Review Task Monitor

THRU:

Harold Brown, Ph.D.

ESAT Deputy Project Officer, EPA

FROM:

Kevin Ludwikoski

ESAT Data Reviewer

THRU:

Ronald A. Ross

ESAT Team Manager

DATE:

May 2, 1991

SUBJECT: Review of inorganic data for Laclede Coal Gas Co.

TID# 07-9103-535 ASSIGNMENT# 772

ICF ACCT# 302-26-535-02 ManTech S.O.# 1073-535

ESAT Document No. ESAT-VII-535-0017

These data were reviewed primarily according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," July 1988 revision with changes given in the Region VII Inorganic Data Review Training Manual and EPA memorandums.

The following comments and attached data sheets are a result of the ESAT review, according to EPA policies, of the following data from the contract laboratory.

CASE NO.: 15978

SITE: Laclede Coal Gas Co.

REVIEWER: Kevin Ludwikoski

LABORATORY: SKINER

METHOD NO.:

CS0390I

EPA ACTIVITY NO.: DSX44

MATRIX: SOIL

TOTAL METALS and CN (SOIL)

SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
MGG593	DSX44403	MGH535	DSX44005
MGH531	DSX44001	MGH536	DSX44110
MGH532	DSX44002	MGH537	DSX44111
MGH533	DSX44003	MGH538	DSX44112
MGH534	DSX44004		

GENERAL

This data review assignment covers <u>NINE SOIL</u> samples analyzed for <u>TOTAL METALS</u> and <u>CYANIDE</u> for case number <u>15978</u>. There were no field blanks, field duplicates or PE samples included with this assignment.

1. Technical Holding Times / Preservation

Technical holding times have not been specified for soils.

2. Initial and Continuing Calibration

All percent recoveries were within control limits.

3. Blanks

Several analytes were detected in the blanks. Corresponding sample results were qualified according to the blank rule using five times the highest blank value. Sample results requiring modification are reported as non-detect on the attached data sheets.

TOTAL METALS and CN (SOIL)

	5 x Highest	
<u>Analyte</u>	Blank (mg/kg)	Oualified Samples
Sb	20	DSX44111
Ag	5	None qualified
Ca	15	None qualified
Cri	7	None qualified
Pb	2.2	None qualified
Mg	32	None qualified
Na	65	None qualified
Zn	9	None qualified

4. ICP Interference Check

Recoveries of solution AB analytes were within control limits. Sb, K, Zn, and Na were found but not present in the ICS solution. All absolute values for these analytes were below the CRDL and IDL; therefore, no data were qualified by the ICP interference check sample.

5. Laboratory Control Standard (LCS)

LCS results were within established control limits.

6. Duplicates

The RPD for Pb exceeded control limits. The sample results that were coded J are listed below.

TOTAL METALS and CN (SOIL)

<u>Analyte</u>	Samples Oualified	Code
Pb	DSX44001, -002, -003, -004, -005, -110 -111, -112, -403, -403S, -928C	J

7. Matrix Spike Sample

Sb and Mn were out of range for matrix spike recovery. The samples that had data qualified are listed below.

TOTAL METALS and CN (SOIL)

Analyte	१ Rec	Sample No.	Code
Sb Mn	41.5 26 0 .8	DSX44928C DSX44001, -002, -003, -004, -005, -110 -111, -112, -403, -403L, -928C	J J

Note: Sb in sample DSX44111 was above CRDL, but had been previously "U" coded by the blank rule.

8. ICP Serial Dilution

The ICP serial dilution results were within control limits for all analytes of interest.

9. Furnace Atomic Absorption

The analytical scheme was followed for Furnace AA analysis. MSA correlation coefficients were not acceptable for Se in samples DSX44002, -003, -004, -110, -112 and -403L. Samples DSX44003 and -004 were below the CRDL and therefore not affected. The other samples were "J" coded accordingly.

10. Summary

One analyte in one sample was qualified by the blank rule, one analyte by the duplicate rule, two by the spike rule and four samples had Se "J" coded due to standard addition outliers.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

ICF Technology Incorporated

ESAT Region VII

NSI Technology Services Corp.

NSI Technology Services

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The Bionetics Corp.

MEMORANDUM

TO:

Larry Marchin, Chemist, CLQA/LABO/ENSV

THRU:

Harold Brown, Ph.D. DPO/LABO/ENSV

FROM:

Peggy Cox, QA/QC Chemist/ESAT/ManTech & &

THRU:

Ronald Ross, Manager/ESAT/ManTech

DATE:

Muh 1 April 35, 1990

SUBJECT: Review of data for LACLEDE COAL GAS

TID#:

07-9103-535

ASSIGNMENT#:

767

ICF ACCT#:

302-26-535-02

NSI S.O.#:

1073-535

ESAT DOCUMENT#: ESAT-VII-535-0018

These data were reviewed according to the "Laboratory Data Validation Functional Guidelines for Evaluating Organic Analyses, " February 1, 1988 revision and the "Laboratory Data Validation Functional Guidelines for Evaluating Pesticides and PCBs."

The following comments and attached data sheets are a result of ManTech Environmental Technology, Inc.'s review of the above mentioned data from the contract laboratory.

CASE/SAS NO.: 15978 LABORATORY:

RECMD METHOD NO.: CS03900

CONTRACT NO.: 68-D0-0158 SITE: LACLEDE COAL GAS

EPA ACTIVITY: DSX44

REVIEWER: P. COX

MATRIX: WATER

SMO SAMPLE NO. EPA SAMPLE NO. SMO SAMPLE NO. EPA SAMPLE NO.

GJ881	DSX44203	GK958	DSX44201
GJ882	DSX44204	GK959	DSX44202
**GJ883	**DSX44205	G X9 60	DSX44301
GJ884	DSX44206	GK961	DSX44301D
GÍ885	DSX44206D	G K 962	DSX44302
GJ886	DSX44207F	G X 963	DSX44304
**GJ887	**DSX44208	GK964	DSX44304D
*GJ888	*DSX44209F	GK965	DSX44900P
GJ889	DSX44303	GJ868	DSX44010

Volatile analysis only.

^{**} Rinsate blanks.

GENERAL

Case 15978 contained 18 water environmental and 18 QC samples analyzed for volatiles and base/neutrals and acids at the low level concentration. Two field blanks, three field duplicates, and one performance evaluation sample were included in this data review assignment. Data review was performed at level 2.

1. TECHNICAL HOLDING TIMES and PRESERVATION

- A. Technical holding times for volatile analyses were exceeded by 3-6 days for aromatic compounds in samples DSX44209F, DSX44303, and DSX44010 and all compounds by 16 days in samples DSX44206RE and DSX44206DRE. Benzene, ethylbenzene, and total xylenes in sample DSX44010 were qualified with a "J" code according to the holding time/preservation rules.
- B. Technical extraction and analysis holding times and required preservation were within quality control limit requirements on all samples for base/neutral and acid analyses.

2. GC/MB TUNING

- A. All volatile GC/MS tunings and mass calibrations were within quality control limit requirements for bromofluorobenzene (BFB).
- B. All base/neutral and acid GC/MS tunings and mass calibration were within quality control limit requirements for decafluorotriphenylphosphine (DFTPP).

3. INITIAL and CONTINUING CALIBRATION

- A. Acetone, 2-butanone, 4-methyl-2-pentanone, and 2-hexanone were outside quality control limit requirements for percent relative standard deviation (greater than 30%) on the initial volatile calibrations. Since all associated data were non-detect or qualified by other rules, no data were qualified based on the initial calibration rules.
- B. Acetone, 2-butanone, trans-1,3-dichloropropene, 4-methyl-2-pentanone, 2-hexanone, and carbon disulfide were outside quality control limit requirements for percent difference (%D) (greater than 25%) on the volatile continuing calibrations. Acetone in sample DSX44908M was qualified with a "J" code based on the continuing calibration rules.

- C. 2,4-dinitrophenol, 4-nitrophenol, pentachlorophenol, butylbenzylphthalate, bis(2-ethylhexyl)phthalate, di-n-octyl phthalate, hexachlorocyclopentadiene, 4,6-dinitro-2-methyl phenol, pyrene, and butylbenzylphthalate were outside quality control limit requirements for percent difference (%D) (greater than 25%) on the base/neutral and acid continuing calibrations. 4-nitrophenol and bis(2-ethylhexyl)phthalate in sample DSX44900P and pyrene in sample DSX44010 were qualified with a "J" by the calibration rules.
- D. All compounds were within quality control limit requirements for average relative response factor and response factor on the volatile and base/neutral and acid initial and continuing calibrations.

4. INTERNAL STANDARD RESPONSE

A. All internal standard response areas were within a factor of 2 when comparing samples and their associated continuing calibration response areas.

5. BLANKS

- A. Six method blanks were analyzed for the volatile fraction with methylene chloride and acetone reported at levels greater than the instrument detection limit (IDL) but less than the contract required quantitation limit (CRQL) (except acetone in sample DSX44908M which was greater than the CRQL). Acetone in samples—DSX44301D and DSX44302 were qualified with a "U" code according to the blank rules.
- B. Two method blanks were analyzed for base/neutrals and acids with phenol and bis(2-ethylhexyl)phthalate reported levels greater than the instrument detection limit (IDL) but less than the contract required quantitation limit (CRQL). Since all associated data were non-detect, no data were qualified by the blank rules.
- C. Two field blanks, DSX44207F and DSX44209F and two rinsate blanks, DSX44205 and DSX44208 were analyzed for volatiles with methylene chloride, chloroform, and toluene reported at levels greater than the instrument detection limit (IDL) but less than the contract required quantitation limit (CRQL). Methylene chloride in sample DSX44900P was qualified with a "U" code according to the blank rules.
- D. One field blank, DSX44207F and two rinsate blanks, DSX44205 and DSX44208 were analyzed for base/neutrals and acids with bis(2-ethylhexyl)phthalate reported at levels greater than the instrument detection limit (IDL) but less than the contract required quantitation limit (CRQL). Since all associated data were non-detect, no data were qualified by the blank rules.

6. SURROGATE RECOVERY

- A. All surrogates were within quality control limit requirements for percent recovery for the volatile fraction except toluene-d8 in samples DSX44206D and DSX44206MS. Upon reanalysis of the samples, the surrogates were within quality control limit requirements. Data was reported from the original sample due to the holding time criterion being exceeded and since all compounds were non-detect, no data were qualified by the surrogate recovery rules.
- B. All base/neutral and acid surrogates were within quality control limit requirements for percent recovery.

7. MATRIX SPIKE/MATRIX SPIKE DUPLICATE

- A. A matrix spike/matrix spike duplicate was analyzed for volatiles and was within quality control limit requirements for percent recovery and relative percent difference.
- B. A matrix spike/matrix spike duplicate was analyzed for base/neutrals and acids with 4-nitrophenol outside quality control limit requirements for percent recovery in the matrix spike and matrix spike duplicate and 2,4-dinitrotoluene for percent recovery in the matrix spike duplicate. All spiking compounds were within quality control limits for relative percent difference. No data were qualified by the matrix spike/matrix spike duplicate rules.

8. PERFORMANCE EVALUATION SAMPLE

- A. Performance evaluation sample, DSX44900P, was submitted to the laboratory for volatile analysis with all compounds contained in the audit identified except dichloromethane and 2-chloroethyl ethyl ether. O-dichlorobenzene was reported as an isomer. No data were qualified by the performance evaluation sample.
- B. Performance evaluation sample, DSX44900P, was submitted to the laboratory for base/neutral and acid analysis with all compounds contained in the audit identified except 4-chlorobenzotrifluoride, m-chlorotoluene, 2,4-dichlorotoluene, 1,3,5-trichlorobenzene or 1,2,3,-trichlorobenzene, and 2,4,6-trichloroaniline. 1,3,5-trichlorobenzene or 1,2,3-trichlorobenzene was reported as an unknown trichlorobenzene and 1,2,4,5-tetrachlorobenzene and 1,2,3,4-tetrachlorobenzene as an unknown tetrachlorobenzenes. Di-n-butylphthalate and bis(2-ethylhexyl)phthalate were also reported. No data were qualified by the performance evaluation sample.

9. COMPOUND IDENTIFICATION and QUARTITATION

- A. All target compound identifications were supported by good agreement between sample and standard mass spectra submitted for each positive sample response.
- B. In a level 2 review, no calculations from raw data (quantitation reports) are performed.
- C. Several compounds in each fraction were "U" coded due to the concentration being greater than the instrument detection limit (IDL) but less than the contract required quantitation limit (CRQL).

10. SUNCARY

- A. Benzene, ethylbenzene, and total xylenes in sample DSX44010 were qualified with a "J" code due to the holding times being exceeded.
- B. Acetone in volatile sample DSX44908M was qualified with a "J" code due to the continuing calibration criterion for percent difference being exceeded.
- C. 4-nitrophenol and bis(2-ethylhexyl)phthalate in base/ neutral and acid sample DSX44900P and pyrene in sample DSX44010 were qualified with a "J" due to the continuing calibration criterion for percent difference being exceeded.
- D. Acetone in samples DSX44301D and DSX44302 and methylene chloride in sample DSX44900P were qualified with a "U" code due to blank contamination.
- E. This data package generally meets the requirements for precision, accuracy, and completeness as described in SOW for Organic Analysis dated March 1990, with the exceptions noted above.

ENVIRONMENTAL SERVICES ASSISTANCE TEAM - ZONE II

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66115

The Bionetics Corp.

TO:

Larry Marchin/Barry Evans

Data Review Task Monitor

Harold Brown, Ph.D. THRU:

NSI Technology Services Corp.

ESAT Deputy Project Officer, EPA

FROM:

Kevin Ludwikoski
ESAT Data Reviewer

Ronald A. Ross THRU:

ESAT Team Manager

DATE: May 7, 1991 SUBJECT: Review of inorganic data for Laclede Coal Gas Co.

TID# 07-9103-535 ASSIGNMENT# 779

ICF ACCT# 302-26-535-02 ManTech S.O.# 1073-535

ESAT Document No. ESAT-VII-535-0026

These data were reviewed primarily according to the "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," July 1988 revision with changes given in the Region VII Inorganic Data Review Training Manual and EPA memorandums.

The following comments and attached data sheets are a result of the ESAT review, according to EPA policies, of the following data from the contract laboratory.

CASE NO.: 15978

LABORATORY: SKINER

SITE: Laclede Coal Gas Co. REVIEWER: Kevin Ludwikoski

METHOD NO.: CS0390I

_ EPA ACTIVITY NO.: DSX44 MATRIX: WATER

TOTAL METALS and CN (WATER)

SMO Sample No.	EPA Sample No.	SMO Sample No.	EPA Sample No.
MGG575	DSX44203	MGG582	DSX44206D
MGG577	DSX44204	MGG584	DSX44207F
MGG579	DSX44205F	MGG586	DSX44208F
MGG580	DSX44206	MGG588	DSX44303
(MGG581	DSX44206) *		

^{*} See discussion

GENERAL

This data review assignment covers <u>EIGHT WATER</u> samples analyzed for <u>TOTAL METALS</u> and <u>CYANIDE</u> for case number <u>15978</u>. There was one field blank (DSX44207F), two rinsates (DSX44205F and DSX44208F), and one field duplicate included with this assignment.

1. Technical Holding Times / Preservation

Technical holding times were observed for all analytes.

2. Initial and Continuing Calibration

All percent recoveries were within control limits.

3. Blanks

Several analytes were detected in the blanks. Corresponding sample results were qualified according to the blank rule using five times the highest blank value. Sample results requiring modification are reported as non-detect on the attached data sheets.

TOTAL METALS and CN (WATER)

	2 x Highest	
<u>Analyte</u>	Blank (ug/L)	Oualified Samples
Al	1550	None qualified
Ag	11	None qualified
Ca	250500	DSX44203, -204, -206, -206L,
		-206D, -208F, -303, -920C
cr	78	DSX44206D, -303
Co	12	None qualified
Cu	985	DSX44203, -204, -206, -206L,
		-206D, -206S, -920C
Pb	158	DSX44205F, -207F, -303, -920C
Mg	96500	DSX44203, -204, -206, -206D,
		-206L, -303, -920C
Na	92500	DSX44204, -303, -920C
N1	43	None qualified
K	19900	DSX44206, -206D, -206L, -303
zn	316	DSX44205F, -207F, -303
As	10	None qualified
Ba	247	None qualified
Fe	24850	DSX44204, -205F, -920C
' Mn	196	DSX44205F
CNT	64	None qualified

4. ICP Interference Check

Recoveries of solution AB analytes were within control limits. K, Sb, Zn, and Na were found but not present in the ICS solution. All absolute values for these analytes were below the CRDL and IDL; therefore, no data were qualified by the ICP interference check sample.

5. Laboratory Control Standard (LCS)

LCS results were within established control limits.

6. <u>Duplicates</u>

The RPD for As exceeded control limits. The sample results that were coded J are listed below.

TOTAL METALS and CN (WATER)

<u>Analyte</u>	Samples Oualified	Code
As	DSX44203, -206, -206D, -206S, -920C	J

7. Matrix Spike Sample

As, Cr, Se, Tl and CN were out of range for matrix spike recovery. The samples that had data qualified are listed below.

TOTAL METALS

<u>Analyte</u>	१ Rec	Sample No.	Code
As	63.4	DSX44203, -206, -206D, -206L, -920C	J
Cr	74.3	DSX44208F -920C	J
Se	0.0	DSX44203, -204, -205, -206, -206D, -206L -207F, -208, -303	I
		DSX44920C	J
Tl	14.5	DSX44203, -204, -205, -206, -206D, -206L -207F, -208, -303	I
		DSX44920C	J
CN	66.9	DSX44203, -204, -205	J

Note: Samples DSX44206D and DSX44303 were above the CRDL, but had been previously "U" coded for Cr according to the blank rule.

Note: All samples that were "J" coded for As, with the exception of DSX44206L, had been "J" coded for As according to the duplicate rule.

8. ICP Serial Dilution

The ICP serial dilution results were within control limits for all analytes of interest.

9. Furnace Atomic Absorption

The analytical scheme was followed for Furnace AA analysis. MSA correlation coefficients were acceptable for As in sample DSX44206 (cc=.9970).

10. Discussion

Sample numbers MGG580 and MGG581 both had an EPA sample number DSX44206. Sample MGG581 had been filtered in the field for Dissolved Metals analysis but was included in the Total Metals package. Both aliquots were analyzed as different samples for Total Metals. Therefore, sample MGG580 was reported as DSX44206 and MGG581 was not reported.

11. Summary

Ten analytes had samples that were qualified by the blank rule, one by the duplicate rule, and five by the spike rule.



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 7 25 FUNSTON ROAD KANSAS CITY, KANSAS 66115

DATE:	3/26/91
SUBJECT:	Activity Number: DSX44 Site Description: Laclada Coal Gas D. Simmons D Simmons CLOA-LARD FNSY
FROM:	D. Simmons CLQA-LABO-ENSY
TO:	LABO Primary File
Com	ments regarding the subject activity are as follows:
•	
#30	01 TPHs sample was
	broken.

```
SAMP. NO
                             = SAMPLE IDENTIFICATION NUMBER
      SAMP NO = SAMPLE IDENTIFICATION NUMBER
QCC = QUALITY CONTROL SAMPLE/AUDIT CODE
M = MEDIA OF SAMPLE (A=AIR, T=TISSUE, H=HAZARDOUS
MATERIAL. S=SEDIMENT/SOIL, W=WATER)
AIRS/STORET LOC. NO. = A SAMPLING SITE LOCATION
IDENTIFICATION NUMBER
BEG. DATE = THE DATE SAMPLING WAS STARTED
BEG. TIME = THE TIME SAMPLING WAS STARTED
END. DATE = THE DATE SAMPLING WAS STOPPED
END. TIME = THE TIME SAMPLING WAS STOPPED
A = PESERVED
       A = RESERVED
      B = RESERVED

B = RESERVED

PES = PESTICIDES BY CONTRACT

= DIOXINS/FURANS BY EPA

E = EXPLOSIVES BY CONTRACT

FLD = FIELD MEASUREMENTS BY EPA
      G = MINERALS & DISSOLVED MATERIALS BY EPA
HER = HERBICIDES BY EPA
     I = IGN CHROMATOGRAPHY ANALYSES BY EPA
MC = METALS BY CONTRACT
BNC = BASE NEUTRALS BY CONTRACT
L = FISH PHYSICAL DATA BY EPA
MET = METALS BY EPA
     MET = METALS BY EPA

N = FISH TISSUE PARAMETERS BY EPA

VC = VOLATILES BY CONTRACT

P = PESTICIDES BY EPA

Q = FLASH POINT ANALYSES BY EPA

R = RESERVED
      BH = SEMIVOLATTIE BY EPA
T = CYANIDE PHENOL BY EFA
      U = RESERVED
      VOA = VOLATILE ORGANICS BY EPA
      HC = HERBICIDES BY CONTRACT
           ⇒ RESERVED
           = RESERVED
      TRK = ACTIVITY TRACKING PARAMETERS BY EPA
     STORET DETECTION IDENTIFIERS
     BLANK NO REMARKS
     J DATA REPORTED BUT HOT VALID BY APPROVED OF PROCEDURES
     I = INVALID SAMPLE/DATA = VALUE NOT REPORTED
     IL IESS THAN (MEASUREMENT DETECTION LIMIT)
    M DETECTED BUT BELOW THE LEVEL FOR ACCURATE QUANTIFICATION
    O PARAMETER HOT AMALIZED
* TOTHERAUTOR/ IN HOUSE / LILLD MEDIA GROUPS
LIELD - + + = AF.HF.SF.TF.WF.ZZ
LONGRACIOR = + = HA.HC.HJ.HK.HO.SC.SJ.SK.SO.SW.TC.TJ.
TK.TO.TW.WA.WG.WE.WJ.WK.WO.WW
IN HOUSE - + = ALL OTHERS
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QUALITY CONTROL AUDIT CODES

A = TRUE VALUE FOR CALIBRATION STANDARD

B = CONCENTRATION RESULTING FROM DUPLICATE LAB SPIKE

C = MEASURED VALUE FOR CALIBRATION STANDARD

D = MEASURED VALUE FOR FIELD DUPLICATE

F = MEASURED VALUE FOR METHOD STANDARD

H = TRUE VALUE FOR METHOD STANDARD

K = CONCENTRATION RESULTING FROM DUPLICATE FIELD SPIKE

L = MEASURED VALUE FOR LAB DUPLICATE

M = MEASURED VALUE FOR LAB BLANK

N = MEASURED VALUE FOR DUPLICATE FIELD SPIKE

P = MEASURED VALUE FOR PERFORMANCE STANDARD

R = CONCENTRATION RESULTING FROM LAB SPIKE

S = MEASURED VALUE FOR LAB SPIKE

T = TRUE VALUE OF PERFORMANCE STANDARD

W = MEASURED VALUE FOR LAB SPIKE

T = TRUE VALUE FOR DUPLICATE LAB SPIKE

Y = MEASURED VALUE FOR FIELD SPIKE

Z = CONCENTRATION RESULTING FROM FIELD SPIKE
                                                    MEDIA CODES
                                                  A = AIR
T = BIOLOGICAL (PLANT & ANIMAL) TISSUE
H = HAZARDOUS MATERIALS/MAN MADE PRODUCTS
                                                   S = SEDIMENT, SLUDGE & SOIL
                                                   UNITS
                                                  NA = NOT APPLICABLE
PG = PICOGRAMS (1 X 10-12 GRAMS)
NG = NANOGRAMS (1 X 10-9 GRAMS)
                                                  UG = MICROGRAMS (1 X 10-6 GRAMS)
MG = MILLIGRAMS (1 X 10-3 GRAMS)
M3 = METER CUBED
MPH = MILES PER HOUR
                                                  SCM = STANDARD (1 ATM, 25 C) CUBIC METER
                                                  KG = KILOGRAM
L = LITER
                                                                   = CENTIGRADE DEGREES
                            SU = STANDARD (PH) UNITS

# = NUMBER

LB = POUNDS

IN = INCHES

M/F = MALE/FEMALE

M2 = SQUARE METER

I.D = SPECIES IDENTIFICATION

GPM = GALLONS PER MINUTE

CFS = CUBIC FEET PER SECOND

MGD = MILLION GALLONS PER DAY

1000G= FLOW. 1000 GALLONS PER LOMPOSITE

UMHOS= CONDUCTIVITY UNITS (1/OHMS)

NTU = TURBIDITY UNITS
                                                 SU = STANDARD (PH) UNITS
                                               NTU = TURBIDITY UNITS
PC/L = PICO (1 x 10-12) CURRIES PER LITER
                                              MV = MILLIVOLT
SQ FT= SQUARE FEET
P/CM2= PICOGRAMS PER SQ. CENTIMETER
U/CM2= MICROGRAMS PER SQ. CENTIMETER
```

ANALYSIS REQUEST REPORT

FOR ACTIVITY: DSX44

SPFD

06/12/91 16:34:32

. FINAL REPORT

FY: 91 ACTIVITY: DSX44

DESCRIPTION: LACLEDE COAL GAS

LOCATION:

MISSOURI

STATUS: ACTIVE

TYPE: SAMPLING - CONTRACT LAB ANALYSIS

PROJECT:

AIRS/

A34

LABO DUE DATE IS 5/14/91. REPORT DUE DATE IS 5/28/91.

INSPECTION DATE: 3/14/91 ALL SAMPLES RECEIVED DATE: 03/15/91

ALL DATA APPROVED BY LABO DATE: 06/12/91

FINAL REPORT TRANSMITTED DATE: 06/21/91

EXPECTED LABO TURNAROUND TIME IS 60 DAYS

EXPECTED REPORT TURNAROUND TIME IS 75 DAYS

ACTUAL LABO TURNAROUND TIME IS 89 DAYS

ACTUAL REPORT TURNAROUND TIME IS 99 DAYS

SAMP.			•	SAMPLE	//	A	CTATE	STORET LOC NO	BEG. Date	BEG. Time	END. Date	END. Time
NO.	QCC	M	DESCRIPTION	STATUS	CONT.	CITY	STATE	FOC NO	D. 1.	,	•	
	400	•••	and a super best to	•	4 ST.	LOUIS	MISSOURI		03/07/91	15:36	1, 1,	
001		S	BACKGROUND SAMPLE-CSI-A	•	4 ST.	Louis	MISSOURI		03/07/91	11:00	/, /,	•
002		5	BACKGROUND SAMPLE-CSI-B		4 St.	Louis	MISSOURI		03/07/91	16:45	/, /,	
003		S	BACKGROUND SAMPLE-CSI-C	- 1	4 St.	Louis	MISSOURI		03/07/91	16:55	1, 1,	•
004		S	BACKGROUND SAMPLE-CSI-D	4	4 ST.	Louis	MISSOURI		03/07/91	17:15	1, 1,	•
005		Ş	BACKGROUND SAMPLE-CSI-E	í	4 St.	LŎŬĬŠ	MISSOURI		03/08/91	09:20	1, 1,	•
006		Ş	BORING LOCAL . W6-B-06-CS2	i	4 St.	Louis	MISSOURI		03/08/91	09:35	', ',	:
007		S	BORING LOCAL #6-R-06-C52	i	4 ST.	Louis	MISSOURI		03/08/91	11:07	4. 4.	:
008		5	R-06-CS3 SAMPLE LOCATION	i	4 St.	ĽŎŬĬŠ	MISSOURI		03/08/91	13:50	', ',	:
009		5	B-14-CS4 SAMPLE LOCATION	i	4 ST.	LÕŪĪŠ	MISSOURI		03/08/91	14:20	', ',	:
010		M	B14-CS4 SAMPLE LOCATION	i	4 ST.	LOUIS	MISSOURI		03/08/91	17:35	', ',	:
011		5	B-07-CS6 SAMPLE LOCATION	i	4 ŠŤ.	LOUIS	MISSOURI		03/09/91	09:24	', ',	•
014			R13-157 SAMPLE LOCATION	ì	4 ŠŤ.	LOUIS	MISSOURI		03/09/91	09:45 13:05	', ',	•
013		÷	BI3 CS7 SAMPLE LOCATION BORING LOCATION-017MOUND STR. POWER	i	3 51	LOUIS	MISSOURI		03/04/91	14 20	1, 1	
101			BORTHO LOCATION 02	i	ã ŠŤ.	LOUIS	MISSOURI		03/02/91	15:30	', ',	
102		٠,٠	BORING LOCALION 02 SURFACE SOLE AT EAST CENTRALOF TANK	i	ã ŠŤ.	LOUIS	MISSOURI		03/04/91	17:25	', ',	
103			BORIFIG FOLATION 05	i	3 51.	LOUIS	MISSOURI		03/04/91	13:50	1, 1,	
104			BETWEEN TANKS 7-8/BORING 07	i	3 ŠŤ.	LOUIS	MISSOURI		03/05/91	14:10	', ',	;
105		},	BORING LOCATION- 15	i	ĬŽ Ě.	LOUIS	MISSOURI		03/06/91	15.15	', ',	:
106		2	BORING LOCATION -16	i	3 ST.	LOUIS	MISSOURI		03/06/91	15:45	1 1	
107		ڔ	BORING LOCATION-17	1	ã ST.	LOUIS	MISSOURI		03/06/91 03/06/91	16:53	1 1	•
108		5	BORING LOCATION-18	1	3 ST.	LOUIS	MISSOURI		03/00/91	07:45	', ',	:
109		Č	R19-SAMPLE LOCATION	1	3 ST.	LOUIS	MISSOURI		03/14/91	07:30	7 7	:
110		۲	R20-SAMPLE LOCATION	1	3 ST.	LOUIS	MISSOURI		03/13/91	17:30	7 7	:
111		5	B21-SAMPLE LOCATION	1	3 ST.	LOUIS .	MISSOURI		00/10/91	, , , , ,	• •	
112		د	DE I DUMI LE FOUTTE DE									

SAMP. NO.	QCC	M		IPLE TUS	CONT.	CITY	STATE	AIRS/ STORET LOC NO	BEG. DATE	BEG. TIME	END. DATE	END. Time
201		W	SAMPLE LOCATION-MULLANPHY STR. /RR TRAC			LOUIS	MISSOURI		03/05/91	17:15	/ /	:
202		W	SAMPLE LOCATION-TERMINAL RR ASSOC. ETC	1		LOUIS	MISSOURI		03/06/91	10:00	/ /	:
203		W	ENGINEERING TEST WELL(147 FT. NORTH)	1		LOUIS	MISSOURI		03/07/91	14:30	1, 1,	:
204 205	_	W	ENGINEERING TEST WELL-363FT.NORTH]		LOUIS	MISSOURI		03/07/91	16:45	/, /,	:
205	r	W	EQUIPMENT RINSATE BLANK	Ţ		roniz	MISSOURI		03/09/91	08:30	/, /,	:
206	_	W	ENGINEERING TEST WELL-20FT. WEST	!		LOUIS	MISSOURI		03/09/91	10:27	/, /,	•
206 207	ñ		ENGINEERING TEST WELL-20FT. WEST	!		roniz	MISSOURI		03/09/91	10:27	/, /,	•
207	ŗ		LACLEDE COAL GAS FIELD BLANK	!		roniz	MISSOURI		03/09/91	12:18	/, /,	•
208 209	בַ	.	GEOPROBE PIPE RINSATE BLANK SAMPLE	!		Louis	MISSOURI		03/08/91	11:00	/, /,	•
209	r		LACLEDE COAL GAS TRIP BLANK	!		roniz	MISSOURI		03/09/91	12:30	1, 1,	•
301			SAMPLE LOCATION-ABANDONED PUMP HOUSE		9 51.	LOUIS	MISSOURI		03/06/91	14:45	· /, /,	•
301	D		DUP. SAMPLE-LOC/ABANDONED PUMP HOUSE		9 SŢ.	LOUIS	MISSOURI		03/06/91	14:45	', ',	•
302 303			SAMPLE LOCATION-UPSTREAM OF NE CORN.	- !		LOUIS	MISSOURI		03/06/91	17:15	', ',	•
303			SW SAMPLE/MISS. RIVER			roniz	MISSOURI		03/07/91	14:00	1, 1,	· ·
304	•		SAMPLE LOCATION-ILL.AMERICAN DRNK.WATR	!		roniz	MISSOURI		03/07/91	08:55	', ',	•
304	D	7	DUP. SAMPLE-LOC/ILL. AMERICAN DRNK.H20	- 1		FORIS	MISSOURI		03/07/91	08:55	', ',	•
401	6	۶	SEDIMENT SAMPLE/GRAB SAMPLE			FORIS	MISSOURI		03/06/91	15:30	<i>'. '.</i>	
401 402	D	5	SEDIMENT SAMPLE/GRAB SAMPLE			LOUIS	MISSOURI	•	03/06/91	15:30	', ',	•
402 403		5	SEDIMENT SAMPLE/GRAB-SW SAMPLE#302			LOUIS	MISSOURI		03/06/91	17:35	', ',	
703 501		5	SEDIMENT SAMPLE/GRAB/MISS RIVER			LOUIS	MISSOURI		03/07/91	14:15 09:15	', ',	:
201		•	IALITUT LUAE GASTUII SAMPIT	u		E CHILLS	M I // II)K I		U.37U9/91	UM: ID	, ,	

ACTIVITY: 1-DSX44

		COMPOUND	UNITS	001	002		003		004		005	5
SMO1	SILVER	BY ICAP	MG/KG 2.4	1 U	2.3	υ	: 2.6	U	2.6	 U	2 9	υ υ
SMO2	ALUMINUM	BY ICAP	MG/KG: 750	00	8600		13000		15000		14000	
• SM03	ARSENIC	BY ICAP	MG/KG: 7.7	,	5.7		8.1		10		:6.7	
• SMO4	BARIUM	BY ICAP	MG/KG: 130)	120		170		120		230	
- SM05	BERYLLIUM	BY ICAP	MG/KG: 1.2	. U	1.2	U	1.3	U	1.3	U	:1.5	U
- SMO6 (CADMIUM	BY ICAP	:MG/KG:1.2	U	1.2	U	1.3	U	1.3	U	:1.5	U
SMO7	COBALT	BY ICAP	MG/KG: 12	V	12	U	13	U	13	U	: 15	U
. SMO8 (CHROMIUM	BY ICAP	MG/KG 11		9.9		14		14		: 15	
SM09 (COPPER	BY ICAP	MG/KG:56		33		26		20		22	
SM10 1	I RON	BY ICAP	MG/KG: 290	00	76000		19000		21000		19000	
SM11 A	MANGANESE	BY ICAP	MG/KG: 320	J	430	J	1700	J	660	J	920	1
SM12 N	MOLYBDENUM	BY ICAP	MG/KG:NA	0	NA	0	: NA	0	NA.	0	NA	0
SM13 N	VICKEL .	BY ICAP	MG/KG: 17		14		20		18		22	
SM14 L	.EAD	BY ICAP	MG/KG: 250	J	78	J	52	J	19	J	26	J
SM15 A	ANTIMONY	BY TCAP	MG/KG: 14	U	14	U	16	U	16	U	18	1)
SM16 S	ELENIUM	BY ICAP	MG/KG:3.0		2.6	J	1.3	U	1.3	U	1 5	U
SM17 T	ITANIUM	BY ICAP	MG/KG NA	0	NA	0	NA	0	NA	0	: NA	0 ,
SM18 T	HALI IUM	BY ICAP	MG/KG: 2.4	U	2.3	U	2.6	U	2.6	U	2 9	U
SM19 V	ANADIUM	BY ICAP	MG/KG: 25		23		33	:	32		.33	
SM20 Z	110	BY TOAP	MG/NG: 120		85	:	65	:	5 3		04	
SMZT C	ALCTUM	B) ICAP	MG/KG:5900	00	47000	: :	16000	:	3200		15000	
5H22 M	AGNESTUM	BY ICAP	:MG/KG:2300)	2500	 -	2900	·	2900	-	1900	
SM23 S	CDTUM	BY ICAP	: MG/KG: 1200) U	1200	U :	1300	U :	1300	· · · ·	1500	u
SM24 P	MULCCATO	BY ICAP	MG/KG: 1200) U	1200	U .	1300	U.	1400		2 100	
5: 01 F	HENOL.		UG/KG: 370	U	400	U :	440	U :	430	U	630	· H
5502 C	ARBA ZOLE		:UG/KG:370		: 400	U :	440	U	430	U	630	U

UG/KG: 370

5528 2.4.6-TRTCHLOROPHENOL

U:440

U:430

U:630

U

VALIDATED DATA

U:400

	COMPOUND	UNITS	001	002	•	003	004	005
5555	BUTYL BENZYL PHTHALATE	UG/KG	370 U	: 400	 U	: 440 U	-: 430 U	:630 U
SS56	3,3'-DICHLOROBENZIDINE	UG/KG	370 U	400	U	440 U	430 U	630 U
SS57	BENZO(A)ANTHRACENE	UG/KG	800	400	U	440 U	:430 U	630 U
SS58	BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG	960 J	1300	J	1200 J	1600 J	2500 J
SS59	CHRYSENE	UG/KG	800	400	U	440	430 U	630 U
SS60	DI-N-OCTYL PHTHALATE	UG/KG	370 U	400	U	440 U	430 U	630 U
SS61	BENZO(B)FLUORANTHENE	UG/KG	1600	400	บ	440 U	430 U	630 U
SS62	BENZO(K)FLUORANTHENE	UG/KG:	400	400	U	440 U	430 U	630 U
5563	BENZO(A)PYRENE	:UG/KG	500	400	U	440 U	430 U	630 U
5564	INDENO(1,2,3-CD)PYRENE	UG/KG	930	400	U	440 U	:430 U	630 11
SS 65	DIBENZO(A,H)ANTHRACENE	UG/KG	370 U	400	U	440 U	430 U	630 U
SS66	BENZO(G.H.I)PERYLENE	UG/KG	200	400	Ū	440 U	430 U	630 U
ST09	CYANIDE	MG/KG:	6	24	3	6.5 U	6.6 U	7.4 U
SV03	CHLOROMETHANE	UG/KG	12 U	12	U	13 U	. 13 U	:14 U
SV04	BROMOME THANE	UG/KG	12 U	: 12	U	13 U	:13 U	:14
SV05	VINYL CHLORIDE	UG/KG:	12 U	12	U	13 U	:13 U	: 14 U
SV06	CHLOROETHANE	UG/KG:	12 U	12	U :	13 U	: 13 U	: 14 U -
SV07	METHYLENE CHLORIDE	: UG/KG: 2	20 U	27	:	13 U	: 13 U	:14 U
SV08	1.1-DICHLOROETHYLENE	: UG/KG: 1	12 U	12	:	13 U	: 13 U	: 14 U
5909	1 1-DICHLOROF THANE	· UG/KG: 1	12 U	:12	: U :	13 "	: 13	14 0
Sv10	TRANS-1, 2-DICHLORUE THYLENE	:UG/KG: 1	2 U	12	: U :	13 U	13 "	14 0 :
5711	CHI OROFORM	UG/KG: 1	2 U	: 12	บ	13 0	່ 13	14 0
SV12	1,2 DICHLOROETHAILE	. UG/KG: 1	2 U	: 12	υ :	13 U	13 U	. 14 0
SVIO	1,1,1-TRICHLORGETHANE	. UG/KG: 1	2 U	. 12	ี บ :	13 U	13 U	.14 ".
SVII	CARBON TETRACHLORIDE	: UG/KG: 1	2 U	: 12	: U :	13 U	13 U	14 U
SV15	BROMOD I CHL OROME THANE	UG/KG: 1	2 U	12	U	13 U	13 U	14 " "

COMPOUND UNITS 001 002 003 004 005 SV16 1.2-DICHLOROPROPANE. UG/KG: 12 U :12 U :13 U :13 SV17 BENZENE : UG/KG: 12 U :12 U :13 U :13 U:14 U : SV18 TRANS-1.3-DICHLOROPROPENE : UG/KG: 12 U :12 U :13 U :13 U:14 U: SV19 TRICHLOROETHYLENE : UG/KG: 12 U :12 U :13 U :13 U:14 U : SV20 CIS-1.3-DICHLOROPROPENE : UG/KG: 12 U : 12 U :13 SV21 DIBROMOCHLOROMETHANE : UG/KG: 12 U:14 U :12 U :13 U :13 SV22 1,1,2-TRICHLOROETHANE : UG/KG: 12 U : 12 U : 13 U :13 U:14 SV24 BROMOFORM : UG/KG: 12 U : 12 U : 13 U :13 SV25 1,1,2,2-TETRACHLOROETHENE UG/KG: 12 U :12 U :13 U :13 U:14 U 13 SV26 TOLUENE : UG/KG: 12 U :12 U :13 11 - 14 SV27 1.1.2.2-TETRACHLOROETHANE : UG/KG: 12 U :12 U :13 U :13 SV28 CHLOROBENZENE : UG/KG: 12 U :12 U :13 U :13 U : 14 U: SV29 ETHYL BENZENE : UG/KG: 12 U :12 U :13 U :13 U:14 **U** : U :53 SV30 ACETONE : UG/KG: 16 U :63 U :40 U :33 J: SV31 CARBON DISULFIDE : UG/KG: 12 U :12 U :13 U · 13 U 14 U SV32 2-BUTANONE : UG/KG: 12 U :12 U :13 U :13 U: U : 14 SV33 VINYL ACETATE : UG/KG: NA O : NA O : NA O:NA O ·NA SV34 2-HEXANONE UG/KG: 12 U : 12 U :13 U : 13 SV35 4-METHYL-2-PENTANONE UG/KG: 12 U : 12 U :13 U :13 U 14 UG/KG: 12 U :12 U :13 0.13SV-16 STYRENE UG/KG: 12 U :12 U : 13 U : 13 SV37 XYLENES, TOTAL 005 2201 SAMPLE HUMBER : NA : 001 : 002 : 003 : 004 2202 ACTIVITY CODE :DSX44 DSX44 NA : DSX44 :DSX44 .DSX44

	COMPOUND	-	DETAIL REI 3-ピイ DO6	PORT AC ジ・パ 007		/: 1-DSX44 デー/ユ 008	√·F′ 009	VALIDATED DATA '.' 'T' O10
SMO1 SILVER	BY ICAP	MG/KG: 2.2	U	2.6	U	2.6 U	J : 2.8	J .
SMO2 ALUMINUM	BY ICAP	MG/KG: 2500		8600		9300	4500	
SMO3 ARSENIC	BY ICAP	MG/KG:4.3		8.5		18	8.5	
SMO4 BARIUM	BY ICAP	MG/KG:92		260		160	250	
SMO5 BERYLLIUM	BY ICAP	:MG/KG:1.1	U	1.3	U	1.3 U	1.4	J
SMO6 CADMIUM	BY ICAP	:MG/KG:1.1	U	1.3	U	1.3 U	1.4	,
SMO7 COBALT	BY ICAP	:MG/KG:11	U	13	Ü	: 13 U	:14 U)
SMO8 CHROMIUM	BY ICAP	:MG/KG:9.5		12		13	: 14	
SMO9 COPPER	BY ICAP	:MG/KG:32		20		27	21	
SM10 IRON	BY ICAP	MG/KG: 7200		17000		41000	16000	
SM11 MANGANESE	BY ICAP	MG/KG:170		720		430	: 180	
SM12 MOLYBDENUM	BY ICAP	MG/KG:NA	0	NA .	0	NA O	NA C	
SM13 NICKEL	BY ICAP	:MG/KG:9.0	U	20		16	11 U	· · · · · · · · · · · · · · · · · · ·
SM14 LEAD	BY ICAP	MG/KG:78	J	:55	J	57 J	580 J	::
SM15 ANTIMONY	BY ICAP	MG/KG: 13	U	16	U	16 U	. 17 U	
SM16 SELENIUM	BY ICAP	:MG/KG: 2.1		1.3	U	1.3 U	1.4 U	
SM17 TITANIUM	BY ICAP	MG/KG: NA	0	NA	0	NA O	:NA O	
SM18 THALLIUM	BY ICAP	:MG/KG:2.2	U	2.6	U	2.6 U	2.8 U	
SM19 VANADIUM	BY ICAP	:MG/KG:11	U	22		29	: 18	
SM20 2100	BY TEAP	MG/KG: 110	J	53	J	71 J	200	
M21 CALCIUM	BY ICAP	:MG/KG: 35000		5200		3100	25000	
SM22 MACHESTUM	BY ICAP	.MG/KG:3100		3000		2400	1800	
SM23 SODIUM	BY ILAP	MG/KG:1100	U	1300	U	1300 U	:1400 U	
SM24 POTASSIUM	BY ICAP	. MG/KG: 1100	U	1300	U	1300 U	1400 U	
SSO1 PHEHOL		UG/KG: 11000	U	12000	U	13000 11	12000 U	
SSO2 CARBAZOLE		:UG/KG:11000	U					
		; ; - 	:					

ACTIVITY: 1-DSX44

COMPOUND	UNITS	006	007	008	009	010
SS29 2,4,5-TRICHLOROPHENOL	UG/KG: 58	3000 U	:59000	บ : 69000 เ	ປ : 59000 U :	
SS30 2-CHLORONAPHTHALENE	UG/KG: 11	000 U	12000	บ 13000 เ	J 12000 U	
SS31 2-NITROANILINE	UG/KG 58	1000 U	59000	บ 69000 เ	59000 U	
SS32 DIMETHYLPHTHALATE	:UG/KG:11	000 Ü	12000	U 13000 , L	1 12000 U	
SS33 ACENAPHTHYLENE	UG/KG:11	000 U	12000	U 15000	: 12000 U	
SS34 3-NITROANILINE	:UG/KG:58	000 U	59000	บ :69000 บ	:59000 U :	
SS35 ACENAPHTHENE	UG/KG:11	000 U	12000	U 32000	12000	
SS36 2.4-DINITROPHENOL	UG/KG:58	000 U	59000	บ :69000 บ	:59000 U :	
SS37 4-NITROPHENOL	UG/KG:58	000 U	59000	N : 68000 N	59000 U	
SS38 DIBENZOFURAN	:UG/KG:11	000 U	12000	U 25000	1200G U	
SS39 2.4-DINITROTOLUENE	UG/KG: 110	000 U	12000	บ 13000 บ	12000 U	
SS40 2.6-DINITROTOLUENE	UG/KG: 110	oóo u	12000	บ : 13000 บ	:12000 U	
SS41 DIETHYLPHTHALATE	UG/KG: 110	000 U	12000	บ : 13000 บ	: 12000 U	
SS42 4-CHLOROPHENYL PHENYL ETHER	UG/KG: 110	000 U	12000	บ :13000 บ	12000 U	
SS43 FLUORENE	UG/KG 200	000	12000	U AND THE REAL PROPERTY.	3000	:
SS44 4-NITROANILINE	UG/KG:580	000 U	59000	บ 69000 บ	:59000 U	
SS45 4.6-DINITRO-2-METHYLPHENOL	UG/KG: 580	000 U	59000	บ 69000 บ	59000 U	•
SS46 N-NITROSODIPHENYLAMINE	UG/KG: 110	000 U	12000	บ 13000 บ	12000 U	
SS47 4-BROMOPHENYL PHENYL ETHER	UG/KG: 110	000 U	12000	บ :13000 บ	:12000 U	
SOUR HEXACHLOROBETIZENE	UG/KG 110	000 U	12000	U 13000 U	: 12000 U	
SAS PENTACHEOROPHENOL	UG/KG:580)00 U	59000 U	U :69000 U	59000 U	
SSEO PHENANTHRENE	.ue/ke: 63 0	000	20000	THE PARTY OF THE P	6000	
SSST ANTHRACENE	UG/KG: 110	000 U	12000 l	The state of the s	12000 U	
SS52 DI-N BUTYL PHTHALATE	UG/KG: 110	00 U	12000 L	J 13000 U	12000 U	
SS53 FLUORANTHENE	UG/KG 110		12000		:12000 U	
SS54 PYRENE	UG/KG:480	00				

ANALYSIS REQUEST DETAIL REPORT ACTIVITY: 1-DSX44 VALIDATED DATA

COMPOUND	UNITS	006	007	008	009	010
SSO3 BIS(2-CHLOROETHYL) ETHER	UG/KG	11000 U	12000 U	: 13000 U	12000 U	:
SSO4 2-CHLOROPHENOL	UG/KG	11000 U	12000 U	13000 U	: 12000 ປ	:
SSO5 1.3-DICHLOROBENZENE	UG/KG	11000 U	12000 U	13000 U	12000 U	
SSO6 1.4-DICHLOROBENZENE	UG/KG:	11000 U	12000 U	:13000 U	12000 U	
SSO7 BENZYL ALCOHOL	UG/KG:	NA O	NA O	NA O	NA O	:
SSO8 1.2-DICHLOROBENZENE	:UG/KG:	11000 U	. 12000 ป	13000 U	12000 U	:
SSO9 2-METHYLPHENOL (O-CRESOL)	:UG/KG:	11000 ປ	: 12000 U	:13000 ປ	12000 U	:
SS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/KG	11000 U	12000 U	: 13000 U	12000 U	
SS11 4-METHYLPHENOL (P-CRESOL)	UG/KG:	11000 U	12000 U	13000 U	12000 U	:
SS12 N-NITROSO-DIPROPYLAMINE	UG/KG:	11000 U	12000 U	13000 U	12000 U	:
SS13 HEXACHLOROETHANE	UG/KG:	11000 U	1.2000 U	13000 U	12000 U	
SS14 NITROBENZENE	UG/KG:	11000 U	12000 U	13000 U	12000 U	:
SS15 ISOPHORONE	UG/KG:	11000 U	12000 U	13000 U	12000 ป	
SS16 2-NITROPHENOL	UG/KG:	11000 U	12000 ป	13000 U	12000 U	
SS17 2 4-DIMETHYLPHENOL	UG/KG:	11000 U	12000 U	: 13000 U	: 12000 U	··
SS18 BENZOIC ACID	UG/KG:	IA O	NA O	NA O	NA O	
SS19 BIS(2-CHLOROETHYOXY) METHANE	UG/KG:1	1000 U	12000 U	13000 U	12000 ປ	•
SS20 2.4-DICHLOROPHENOL	UG/KG: 1	1000 U	12000 ป	: 13000 U	12000 U	
SS21 1.2.4-TRICHLOROBENZENE	UG/KG: 1	1000 U	12000 U	13000 U	12000 U	
SSZZ NAPHTHALENE	UG/KG	3000	87000	2200000	68000	
5523 4-CHEUROANTEINE	UG/KG: 1	1000 U	12000 U	13000 U	12000 U	
SS24 HEXACHI OROBUTADI ENE	UG/KG: 1	1000 U	12000 ປ	: 13000 U	12000 U	·
SS25 4 CHLORO~3-METHYLPHENOL	UG/KG: 1	1000 U	12000 U	13000 ป	12000 U	
SS26 2-METHYLNAPHTHALENE	UG/KG	0000	43000	69000	84000	
SS27 HEXACHLOROCYCLOPENTADIENE	UG/KG: 1	1000 U	12000 U	13000 U	120G0 II	
SS28 2.4.6-TRICHLOROPHENOL	UG/KG: 1	1000 U	12000 U	13000 U	12000 U	
	::-	~~~~~~~~			;	

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ACTIVITY: 1-DSX44

COMPOUND	UNITS	006	007		008	009	010
SS55 BUTYL BENZYL PHTHALATE	UG/KG:	11000 Ų	12000	 υ	: 13000 U	12000	U :
SS56 3,3'-DICHLOROBENZIDINE	UG/KG:	11000 U	12000	U	13000 U	12000	U
SSS7 BENZO(A)ANTHRACENE	UG/KG:	18000	12000	U	120000 1 120000	12000	U
SS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG	11000 U	12000	v	13000 U	12000	U
SS59 CHRYSENE	UG/KG:	28000	12000	U	110000167025	16000	
SS60 DI-N-OCTYL PHTHALATE	UG/KG:	11000 U	12000	U	13000 U	12000	U
SS61 BENZO(B)FLUORANTHENE	UG/KG:	11000 ປ	12000	U	61000	12000	U
SS62 BENZO(K)FLUORANTHENE	UG/KG	11000 U	12000	V	52000	12000	U :
SS63 BENZO(A)PYRENE	UG/KG	11000 , U	12000	U	95000	12000	U
SS64 INDENOCT 2 3-CD PYRENE	UG/KG	11000 U	12000	U	37000	12000	U
SS65 DIBENZO(A,H)ANTHRACENE	UG/KG:	11000 U	12000	U	13000 U	12000	υ :
SS66 BENZO(G.H.1)PERYLENE	UG/KG: 1	11000 U	12000	U	14000	12000	Ü
STO9 CYANIDE	MG/KG	37	180		860	7.0	U :
SVO3 CHLOROMETHANE	UG/KG: 1	1500 U	12000	U	16000 ป	1500	U :
SVO4 BROMOMETHANE	UG/KG: 1	1500 U	12000	บ	16000 U	1500	U :
SVOS VINYL CHLORIDE	UG/KG: 1	500 U	12000	U	16000 U	1500	U
SVO6 CHLOROETHANE	UG/KG: 1	500 U	12000	U	16000 U	1500	U
SVO7 METHYLENE CHLORIDE	UG/KG: 1	500 U	12000	U:	16000 U	1500	U :
SVO8 1.1-DICHLOROFTHYLENE	UG/KG:1	500 U	12000	U	16000 U	1500	0
was to latenakotawa	UG/KG: 1	500 U :	12000	· ()	16000 U	1500	u :
SVIO TRANS-1, 2-DE HEOROL TOYLENE	UG/KG:1	500 U	12000	U :	16000 U	1500	H ·
SV11 CHLOROFORM	: UG/kG: 1	500 U	12000	U :	16000 11	1500	IJ
SV12 1.2 DICHLORGETHANE	. UG/KG: 1	500 U	12000	U :	1600Q U :	1500	Ū
SV13 1.1.1 TRICHLORGETHANE	UG/KG: 1	500 U	12000	U :	16000 U	1500	Ü .
SV14 CARBON TETRACHLORIDE	UG/KG: 1	500 U :	12000	u :	16000 11	1500	Ū
SV15 BROMODICHLOROMETHANE	UG/KG: 1	500 U	12000	U	16000 U : 1	1500	Ü :

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ACTIVITY: 1-DSX44

	COMPOUND	UNITS	006	007		800	009	010
SV16 1.2-DICHLOR	OPROPANE	UG/KG 1	1500 U	12000	v	16000 ປ	1500 ເ) :
SV17 BENZENE		UG/KG	3500	140000		150000	19000	3
SV18 TRANS-1,3-D	J CHLOROPROPENE	UG/KG: 1	1500 U	12000	U	16000 U	1500 U	
SV19 TRICHLOROET	HYLENE	UG/KG: 1	500 ป	12000	U	16000 U	1500 U	:
SV20 CIS-1.3-DIC	HLOROPROPENE	UG/KG: 1	500 <u> </u>	12000	U	16000 U	1500 U	:
SV21 DIBROMOCHLO	ROME THANE	UG/KG: 1	500 U	12000	U	16000 U	: 1500 U	:
SV22 1.1.2-TRICHL	OROETHANE	UG/KG: 1	500 U	12000	U	16000 U	: 1500 ป	:
SV24 BROMOFORM		UG/KG: 1	500 U	12000	IJ	16000 U	: 1500 U	- ;
SV25 1.1.2.2-TETE	RACHLOROE THENE	UG/KG: 1	500 U	12000	U	16000 U	1500 U	:
SV26 TOLUENE		UG/KG	700	26000		58000	ູ່ 1500 ບ	:
SV27 1.1.2.2-TETR	ACHLOROE THANE	UG/KG 1	500 U	12000	U	16000 U	: 1500 U	:
SV28 CHLOROBENZEN	iE	UG/KG: 1	500 U	12000	Ü	16000 U	1500 U	:
SV29 ETHYL BENZEN	E	UG/KG:2	4000	46000		96000	5500	:
SV30 ACETONE		UG/KG: 1	500 U	12000	U	16000 U	1500 U	
SV31 CARBON DISUL	FIDE	UG/KG: 1	500 U	12000	U	16000 U	1500 1/	:
SV32 2-BUTANONE		UG/KG: 11	500 U	12000	ับ	16000 U	1500 ບ	:
SV33 VINYL ACETAT	E	UG/KG N/	A 0	NA	0	NA O	NA O	:
SV34 2-HEXANONE		UG/KG 19	500 U	12000	U	16000 U	1500 ป	:
SV35 4-METHYL-2-P	ENTANONE	UG/KG 19	500 U	12000	U	16000 U	1500 U	:
SVA6 STYRENE		: UG/KG: 15	500 U	42000	:	16000 11	1500 U	,
SV37 AYLENES, TOTA	AL	UG/KG (2 8	3000	230000			9400	
WMO1 SILVER	BY TOAP	UG/L		*	; :			ATTACHMENT
WMO2 ALUMINUM	BY ICAP	UG/L	:		·: :			ATTACHMENT
WMO3 ARGENIC	BY ICAP	UG/L	:		:			ATTACHMENT
WMO4 BARIUM	BY ICAP	UG/I.	:		:			ATTACHMENT
WMO5 BERYLLIUM	BY ICAP	UG/L :			: :			ATTACHMENT

COMPOUND	UNITS	006	007	008	009	010	
BY ICAP	: UG/L :	: :			:	. ATTACHME	NT
BY ICAP	UG/L	:			:	ATTACHME	NT
BY ICAP	UG/L	:			:	ATTACHME	NT
BY ICAP	UG/L :		:		:	ATTACHME	NT
BY TCAP	UG/L	:	:		:	ATTACHME	NT
E BY ICAP	UG/L :	:	:		:	ATTACHMEN	11
UM BY ICAP	UG/L	::	:	,	:	ATTACHMEN	T
BY ICAP	UG/L .	:	:		:	ATTACHMEN	NT
BY ICAP	UG/L	: :		J		ATTACHMEN	VT
BY ICAP	: UG/L :		:		: 	ATTAL HIME	11
BY ICAP	:UG/L	:	:		:	ATTACHMEN	11
BY ICAP	UG/L	:	:			ATTACHMEN	IT
BY ICAP	UG/L	:	:			ATTACHMEN	IT
BY ICAP ;	UG/L					ATTACHMEN	iT
BY ICAP	UG/L	:	:	:		ATTACHMEN	1
TOTAL BY ICAP	MG/L	·:	:	;		ATTACHMEN	IT.
A, TOTAL BY ICAP	MG/L	:	:: :			ATTACHMEN	1
OTAL BY ICAP	MG/L :		·	:		ATTACHMEN	ī
I. TOTAL BY ICAP	MG/L	:	:			ATTACHMEN	ī
	UG/L :	:	·:-			150	łı
OROETHYL) ETHER	UG/L :	:	· · · · · · · · · · · · · · · · · · ·	:	• • • • • • • • • • • • • • • • • • • •	150	
HENOL	UG/L :	:	:			150	0
OROBENZEHE	UG/L :	······································	:			150	ü
OROBENZENE	UG/L :	:-	·::-			150	Ü
COHOL	UG/L		·	:		: 150	Ü
OROBENZENE	UG/L	:::-:				150	Ü
	BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP E BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP BY ICAP	BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L E BY ICAP UG/L UM BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L TOTAL BY ICAP UG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L COROBENZENE UG/L COROBENZENE UG/L COHOOL UG/L COHOOL UG/L	BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L E BY ICAP UG/L OTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L OROBENZENE UG/L OROBENZENE UG/L OROBENZENE UG/L OROBENZENE UG/L	BY ICAP	BY ICAP UG/L ITOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. TOTAL BY ICAP MG/L A. OROBENZENE UG/L OROBENZENE UG/L COHOL UG/L	BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L BY ICAP UG/L E BY ICAP UG/L E BY ICAP UG/L BY ICAP UG/L	BY ICAP UG/L ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN ATTACHMEN UG/L UG/L

•	COMPOUND	UNITS	006	007	008	009	010)
WSO9 2-METHYL	PHENOL (O-CRESOL)	: UG/L		: : :		:	: 150	u
WS10 BIS(2-CH	LOROISOPROPYL) ETHER	UG/L				:	150	U
WS11 4-METHYL	PHENOL (P-CRESOL)	UG/L		:		:	150	U
WS12 N-NITROS	O-DIPROPYLAMINE	UG/L				:	150	U
WS13 HEXACHLO	ROETHANE	UG/L		,		:	150	Ü
WS14 NITROBEN	ZENE	UG/L				:	150	U
WS15 ISOPHORON	VE	UG/L					150	U
WS16 2-NITROPH	lENOL	UG/L				:	: 150	U
WS17 2.4-DIMET	THYL PHÊNOL	UG/L				:	150	U
WS18 BENZOIC, A	CID	UG/L				:	150	U
WS19 BIS(2-CHL	OROETHYOXY) METHANE	UG/L				:	150	บ
WS20 2.4-DICHL	OROPHENOL	UG/L				·	: 150	U
WS21 1.2.4-TRI	CHLOROBENZENE	UG/L					: 150	: :
WS22 NAPHTHALE	NE ,	UG/L					890	
WS23 4-CHLOROA	NIL INE	UG/L	:	:			:150	ļ)
WS24 HEXACHLOR	OBUTADIENE	UG/L :	: :			:	: 150	Ü
WS25 4-CHLORO-	3-ME THYL PHENOL	UG/L	:	:			150	-U
WS26 2-METHYLN	APHTHALENE	UG/L	:	: :			9 80	
WS27 HEXACHLOR	OCYCLOPENTADIENE	UG/L	:	:			150	U
Wide 2 d 6 (R)	्स जंब (त वर्ष	0671	:	:	:		150	H
W 29 2 4 5-1RD	HLOROPH NO	0671	:	<u></u>	·		150	- 11
WS30 2-CHLOPON	APHTHALLIE	:UG/L	:	:			. 150	11
WL31 2 HEIROAN	ILINE (GRING NITROANILINE)	UG/L	:	;	;		150	υ.
WSG2 DIMETHYLPI	HTHALATE	UG/L	:	:			. 150	U.
WS33 ACEHAPHTH	VI EMC	UG/L :		:			150	H
W534 3-NIIROAN	II. INE	UG/L					150	IJ

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ACTIVITY: 1-DSX44

COMPOUND	UNITS	006	007	008	009	010)
WS35 ACENAPHTHENE	UG/L		·:	:	-;	150	·
WS36 2.4-DINITROPHENOL	UG/L		:	:	-:	150	Ú
WS37 4-NITROPHENOL	UG/L		:	:	:	150	U
WS38 DIBENZOFURAN	UG/L		:	:	:	150	U
WS39 2,4-DINITROTOLUENE	UG/L		:	:	:	150	U
WS40 2.6-DINITROTOLUENE	UG/L		:	:	:	: 150	U
WS41 DIETHYLPHTHALATE	UG/L		:	:	:	: 150	U
WS42 4-CHLOROPHENYL PHENYL ETHER	UG/L		:	:	: :	: 150	U
WS43 FLUORENE	UG/L		:	:	:	170	
WS44 4-NITROANILINE	UG/L :		:	:	:	380	u
WS45 4.6-DINITRO-2-METHYLPHENOL	UG/L		:	:	:	380	U
WS46 N-NITROSODIPHENYLAMINE	UG/L		:	:	:	150	U
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L			:		150	U
WS48 HEXACHLOROBENZENE	UG/L			:	:	150	U
WS49 PENTACHI OROPHENOL	UG/L				:	: 380	U.
WS50 PHENANTHRENE	UG/L			:	:	§ 10	
WS51 ANTHRACENE	UG/L			:	:	150	U
WS52 DI-N-BUTYL PHTHALATE	UG/L :			:	:	150	b
WS53 FLUORANTHENE	UG/L				:	150	U
WS54 PYRENE	UG/L :			:	:	ADS-10	2
WS55 BUTYL BENZYL PHIHALATE	U6/L	:			, · · · · · · · · · · · · · · · · · · ·	150	Ü
WS56 3.37-DICHLOROBENZIDINE	UG/L	:			;	150	U
W35.7 BUNL'O(A)ANTHRACENE	UG/L				. 	150	Ü
WS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/L					. 150	U
W359 CHRYSLNE	UG/L					180	1.
WS60 DI-N-OCTYL PHTHALATE	UG/L	:				150	i

COMPOUND	UNITS	006	007	800	009	010)
WS61 BENZO(B)FLUORANTHENE	:UG/L :		:	· ;		150	U
WS62 BENZO(K)FLUORANTHENE	UG/L		•	:		: 150	U
WS63 BENZO(A)PYRENE	:UG/L		:	:		. 150	U
WS64 INDENO(1,2,3-CD)PYRENE	UG/L		:	:		: 150	U
WS65 DIBENZO(A.H)ANTHRACENE	UG/L		:	:		150	U
WS66 BENZO(G.H.I)PERYLENE	UG/L		:			150	U
WS67 CARBAZOLE	UG/L		:	::		150	U
WT09 CYANIDE, TOTAL	MG/L		:			ATTACHME	NT
WVO3 CHLOROMETHANE	UG/L			:		200	U
WVO4 BROMOMETHANE	UG/L			:		∠00	;;
WVO5 VINYL CHLORIDE	UG/L			:		200	Ü
WVO6 CHLOROETHANE	UG/L			:		200	U
WVO7 METHYLENE CHLORIDE	UG/L					200	U
WVO8 1.1-DICHLOROETHENE	UG/L :					200	U
WVO9 1.1-DICHLOROETHANE	UG/L					200	.0
WV10 1.2-DICHLOROETHENE, TOTAL	UG/L					200	U :
WV11 CHLOROFORM	UG/L					200	. U
WV12 1.2-DICHLOROETHANE	UG/L					200	U :
WV13 1.1.1-TRICHLOROETHANE	UG/L					200	IJ
WV14 CARBON TETRACHLORIDE	UG/I					z00	II.
WV15 BROWODTO HEOROME DIAME	:UG/L	:				حال	(1
WV16 1.2 DICHLOROPROPANE	UG/L	:				200	U
WV17 BLNZENE	UG/L					3100	j
WV19 TRICHLOROETHENE	UG/L	:				200	U
WVZO CIS 1.3-DICHEOROFROPENE	UG/I.	:				رن.	U
WV21 DIBROMOCHLOROMETHANE	UG/I.			:		200	U
		; •					•

COMPOUND	UNIT	s 006	007	008	009	010	•
WV22 1,1.2-TRICHLOROETHANE	UG/L	-: :				200	U
WV24 BROMOFORM	UG/L		:		:	200	บ
WV25 TETRACHLOROETHENE	UG/L	:	:	:	:	200	U
WV26 TOLUENE	UG/L	-: :	:	<u>:</u>	:	200	U
WV27 1.1.2.2-TETRACHLOROETHANE	UG/L	:			:	200	U
WV28 CHLOROBENZENE	UG/L				:	200	Ū
WV29 ETHYL BENZENE	: UG/L	:		:	;	: 280	J
WV30 ACETONE	: UG/L	:	:	:	:	200	U .
WV31 CARBON DISULFIDE	UG/L	:			:	200	U
WV32 2-BUTANONE	UG/L	:	:	:	:	200	U
WV33 VINYL ACETATE	UG/L	:	:	:	:	200	U
WV34 2-HEXANONE	UG/L	:	:	:	:	200	U
WV35 4-METHYL-2-PENTANONE	UG/L	:	:	::	:	200	U :
WV36 STYRENE	UG/L	:	: :	:	:	200	Ū
WV37 XYLENES. TOTAL	:UG/L	:	:	:	:	290	
WV40 TRANS-1.3-DICHLOROPROPENE	:UG/L	:		:	:	200	U
2201 SAMPLE NUMBER	: NA	:006	007	:008	: 009	:010	: :
2ZO2 ACTIVITY CODE	NA NA	:DSX44	: DSX44	DSX44	DSX44	DSX44	: -:

	COMPOUND	UNITS	011	012		013		101		102	
SMO1 SILVER	BY ICAP	MG/KG: 2.	5 U	2.8		3.0	U	2.4	U	2 5	U
SMO2 ALUMINUM	BY ICAP	MG/KG: 71	00	4600		3400		8400		4400	
SMO3 ARSENIC	BY ICAP	MG/KG: 7.	6	12		10		21		12	
SMO4 BARIUM	BY ICAP	MG/KG: 23	0	56		61	U	220		78	
SMO5 BERYLLIUM	BY ICAP	:MG/KG:1.:	3 U	1.4	U	1.5	U	1.2	U	1.3	U
SMO6 CADMIUM	BY ICAP	:MG/KG:1.3	3 · U	1.4	U	1.5	U	1.2	U	1.3	U
SMO7 COBALT	BY ICAP	MG/KG:13	Ū	14	U	15	U	12	U	13	U
SMO8 CHROMIUM	BY ICAP	MG/KG:11		21		17		19		9.6	
SMO9 COPPER	BY ICAP	MG/KG:24		83		: 130		140		36	
SM10 IRON	BY ICAP	MG/KG: 300	000	120000	- -	19000		: 45000		23000	
SM11 MANGANESE	BY ICAP	MG/KG: 260)	210		150		: 350		140	::
SM12 MOLYBDENUM	BY ICAP	MG/KG:NA	0	NA	0	NA	0	NA .	0	NA	0
SM13 NICKEL	BY ICAP	MG/KG:17		24		15		28		17	
SM14 LEAD	BY ICAP	MG/KG: 130	J	160	J	250	J	300	J	120	J
SM15 ANTIMONY	BY ICAP	MG/KG:15	U	17	U	18	U	: 15	U	15	()
SM16 SELENIUM	BY ICAP	:MG/KG:1.6		1.4	U	1.5	U	2.4		3.2	:
SM17 TITANIUM	BY ICAP	MG/KG:NA	0	NA	0	NA	0	NA	0	NA	٠0:
SM18 THALLIUM	BY ICAP	MG/KG 2.5	U	2.8	U	3.0	U	2.0	U	2.5	U
SM19 VANADIUM	BY ICAP	MG/KG:19		24		18		29		18	
SM2O ZINC	BY TEAP	:MG/KG:72	J	79	J	98	.1	700	.1	130	.1
-M21 CALLEUM	B. ILAP	MG/KG: 190	00	19000	:	6200		11000		12000	
SM22 MAGNESTUM	BY ICAP	MG/KG: 180	0	1400	U	1500	U	2000	:	1300	
SM23 SODIUM	BY ICAP	:MG/KG:130	O U	1400	Ū:	1500	U :	1500	:	1300	υ.
SM24 POTASSIUM	BY ICAP	MG/KG: 130	O U	1400	U :	1500	U	1200	U .	1300	:
SSO1 PHENOL		UG/KG: 1300	00 U	13000	U	14000	IJ	12000	U :	14000	0
SSO2 CARBAZOLE		UG/KG:				14000	IJ	12000	U	14000	U :
		: :			:						

COMPOUND	UNITS	011		012		013		101		102	
SSO3 BIS(2-CHLOROETHYL) ETHER	UG/KG	13000	U	13000	U	14000	U	12000	U	14000	U
SSO4 2-CHLOROPHENOL	UG/KG	13000	U	13000	U	14000	U	12000	U	14000	U
SSOS 1,3-DICHLOROBENZENE	UG/KG	13000	U	13000	Ü	14000	U	12000	U	14000	U
SSO6 1,4-DICHLOROBENZENE	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SSO7 BENZYL ALCOHOL	UG/KG	NA	0	: NA	0	NA	0	: NA	0	NA	Ö
SSOB 1.2-DICHLOROBENZENE	UG/KG	13000	U	13000	U	14000	U	12000	U	14000	U
SSO9 2-METHYLPHENOL (O-CRESOL)	UG/KG	13000	U	13000	υ	14000	U	12000	U	14000	U
SS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/KG	13000	U	13000	U	14000	U	12000	U	. 14000	U
SS11 4-METHYLPHENOL (P-CRESOL)	UG/KG	13000	U	13000	U	14000	U	12000	U	14000	U
SS12 N-NITRUSO-DIPRUPYLAMINE	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SS13 HEXACHLOROETHANE	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SS14 NITROBENZENE	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SS15 ISOPHORONE	:UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SS16 2-NITROPHENOL	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	U
SS17 2.4 DIMETHYLPHENOL	:UG/KG:	13000	U	13000	U	14000	U	12000	IJ	14000	ĮĮ
SS18 BENZOIC ACID	UG/KG:I	NA	0	NA .	0	: NA	0	NA	0	NA	0
SS19 BIS(2-CHLOROETHYOXY) METHANE	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	- U
SS20 2.4 DICHLOROPHENGL	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	· U
SS21 1.2.4-TRICHLOROBENZENE	UG/KG	13000	U	13000	U	14000	U	12000	U	14000	U
SSZZ NAPHTHALENE	UG/KG:	80000	· · ·	13000	U	5 5000		12000	IJ	14000	U
\$523 dechioroanti fin	UG/KG:	13000	U	13000	U	14000	U	12000	U	14000	Ü :
SG21 HEXACHI OROBUTADITHE	UG/KG: 1	13000	U	13000	U	: 14000	U	12000	U.	14000	Ü,
5525 4 CHLORO 3: METHYLPHLHOU	UG/KG: 1	3000	U	13000	Ū	14000	U	12000	U :	14000	Ü.
S326 2 METHYLNAPHTHALEHE	.UG/KG 🛂	3000	·	13000	U	14000	U .	12000	U :	1-1000	· · ·
SS27 HEXACHLOROCYCLOPENTADIENE	.UG/KG: 1	3000	U	13000	Ü	14000	U	12000	IJ	14000	U .
5528 2.4.6-TRICHLOROPHENOL	UG/KG: 1	3000	U	13000	Ū	14000	U	12000	U	14000	U

COMPOUND	UNITS	011	012		013	101	102
SS29 2.4.5-TRICHLOROPHENOL	: UG/KG	:66000 I	66000	 U	74000 U	:62000	U 73000 U
SS30 2-CHLORONAPHTHALENE	UG/KG	13000	13000	U	14000 U	12000	U 14000 U
SS31 2-NITROANILINE	UG/KG	66000 L	66000	บ	74000 U	62000	บ 73000 บ
SS32 DIMETHYLPHTHALATE	UG/KG	13000	13000	U	14000 U	12000	U : 14000 U
SS33 ACENAPHTHYLENE	UG/KG	20000	13000	· U	14000 U	12000	U : 14000 U
SS34 3-NITROANILINE	UG/KG	66000 U	66000	U	74000 U	62000	บ :73000 บ
SS35 ACENAPHTHENE	UG/KG	15000	13000	U	22000	12000	บ :14000 บ :
SS36 2.4-DINITROPHENOL	UG/KG	66000 U	66000	U	74000 U	62000	บ : 73000 บ :
SS37 4-NITROPHENOL	UG/KG	66000 U	:66000	บ	74000 U	62000	บ 73000 U
SS38 DIBENZOFURAN	UG/KG	13000 U	13000	U	14000 U	12000	J 14000 U
SS39 2.4-DINITROTOLUENE	UG/KG	13000 U	13000	Ū	14000 U	12000	J :14000 U
SS40 2.6-DINITROTOLUENE	UG/KG	13000 U	13000	U	14000 U	12000	J : 14000 U
SS41 DIETHYLPHTHALATE	UG/KG	13000 U	13000	U	14000 U	12000	J :14000 U
SS42 4-CHLOROPHENYL PHENYL ETHER	UG/KG	13000 U	13000	U	14000 U	12000 U	1 14000 U
SS43 FLUORENE	UG/KG	24000	i: 13000	U	21000	12000 (I 14000 .U
SS44 4-NITROANILINE	UG/KG:	66000 U	66000	Ü	74000 U	62000 U	73000 U
SS45 4.6-DINITRO-2-METHYLPHENOL	UG/KG:	66000 U	66000	U	74000 U	62000 U	73000
SS46 N-NITROSODIPHENYLAMINE	UG/KG:	13000 U	13000	U	14000 U	12000 U	1:14000 U
SS47 4-BROMOPHENYL PHENYL ETHER	UG/KG:	13000 U	13000	: U :	14000 U	12000	14000 U
SS48 HEXACHLOROBENZEME	UG/KG:	13000 U	: 13000	:	14000 0	12000 0	14000 U
5549 PENTACHLOROPHENOL	UG/KG:	66000 U	66000	U	74000 u	62000 0	/3000 U
SSEO PHENANTHRENE	:UG/KG:	& 000	13000	U:	59000	12000 0	14000 U
SSST ANTHRACENE	UG/KG:	3000	: 13000		4	12000 U	14000 U
SS52 DI-H-BUTYL PHTHALATE	UG/KG.	13000 U	13000			12000 U	1 1000 U .
SSSS FI UORANTHENE	:UG/KG:	80000	13000	Ū:	9000: 1.1.1.1	12000 !!	1.4C00 U
SS54 PYRENE	:UG/KG:	210000 J	: 13000	U	5000	12000 0	21000

COMPOUND	UNITS	011	012		013	101	102
SS55 BUTYL BENZYL PHTHALATE	: UG/KG:	13000 U	13000	 U	: 14000 U	: 12000 U	14000 U
SS56 3.3'-DICHLOROBENZIDINE	UG/KG:	13000 U	13000	U	14000 U	12000 U	:14000 U
SS57 BENZO(A)ANTHRACENE	:UG/KG	,99000	13000	Ų	21000	12000 U	:14000 U
SS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG	13000 U	13000	U	14000 U	12000 U	14000 U
SS59 CHRYSENE	UG/KG	94000	13000	U	£1000	:12000 U	:14000 U
SS60 DI-N-OCTYL PHTHALATE	UG/KG:	13000 U	13000	U	14000 U	:12000 U	:14000 U
SS61 BENZO(B)FLUORANTHENE	UG/KG:	76000	: 13000	U	15000	12000 U	14000 U
SS62 BENZO(K)FLUORANTHENE	UG/KG:	89000	: 13000	U	7000	12000 U	14000 U
SS63 BENZO(A)PYRENE	UG/KG:	20000	13000	U	≩ 3000	12000 U	:14000 U
SS64 INDENO(1.2.3-CD)PYRENE	UG/KG:	75000	13000	U	14000 U	12000 ()	14000 11
SS65 DIBENZO(A, H)ANTHRACENE	:UG/KG	13000 U	13000	U	14000 U	:12000 ປ	:14000 U
SS66 BENZO(G.H.I)PERYLENE	:UG/KG	88000	13000	U	14000 U	12000 U	: 14000 U :
STO9 CYANIDE	:MG/KG	?50 ,	170		7.6 U	<u>.</u> 33	.6.3 U :
SVO3 CHLOROMETHANE	UG/KG	3100 ย	: 13	U	1700 U	:	::
SVO4 BROMOMETHANE	UG/KG:	3100 U	: 13	U	1700 U	: :	
SVO5 VINYL CHLORIDE	:UG/KG:3	3100 U	:13	U	1700 U	:	:
SVO6 CHLOROETHANE	UG/KG:	3100 U	:13	U	1700 U	: 	: :
SVO7 METHYLENE CHLORIDE	UG/KG:	3100 U	: 13	U	1700 U	:	
SVO8 1 1-DICHLOROETHYLENE	UG/KG:3	3100 U	: 13	U	1700 U		· ·
SVO9 1 1-DTCHLOROETHANE	UG/KG 3	3100 U	: 13	U	1700 U		
SWTO TRANS-1 2-DTCHLOROETHYLENE	UG/KG:3	3100 U	: 13	U	1700 U	:	:
SV11 CHLOROFORM	UG/KG:3	3100 U	: 13	U	1700 U	:	
SV12 1.2 DICHLOROETHANE	UG/KG 3	3100 U	13	U	1700 U		
SV13 1.1.1-TRICHLOROETHANE	UG/KG 3	3100 U	13	U .	1700 U		
SV14 CARBON TETRACHLORIDE	UG/KG: 3	3100 U	13	U :	1700 U	:	
SV15 BROMODICHLOROMETHANE	UG/KG:3	1100 ປ	13	U	1700 U	:	
	,	·				•	•

COMPOUND	UNITS	01	1	(012	013		101	.102	2
SV16 1.2-DICHLOROPROPANE	. UG/KG	3100	U	13	U	:1700	U	:		
SV17 BENZENE	: UG/KG	48000		13	v	1800		:	:	
SV18 TRANS-1.3-DICHLOROPROPENE	:UG/KG	3100	Ü	:13	U	1700	U		:	
SV19 TRICHLOROETHYLENE	UG/KG	3100	U	13	U	1700	U	:		
SV20 CIS-1,3-DICHLOROPROPENE	: UG/KG	3100	U	13	U	1700	U	:		
SV21 DIBROMOCHLOROMETHANE	UG/KG	3100	Ū	13	U	1700	U	:		
SV22 1.1.2-TRICHLOROETHANE	UG/KG	3100	U	13	U	1700	U	:		
SV24 BROMOFORM	UG/KG	3100	U	13	U	1700	U	:	:	
SV25 1.1.2.2-TETRACHLOROETHENE	UG/KG	3100	U	13	U	1700	U	;		
SV26 TOLUENE	UG/KG	3100	U	13	U	1700	U	;	:	
SV27 1.1.2.2-TETRACHLOROETHANE	: UG/KG	3100	U	13	U	1700	U	:	:	
SV28 CHLOROBENZENE	UG/KG	3100	U	13	ย	1700	U	:		
SV29 ETHYL BENZENE	UG/KG	34000		13	U	1700	U			
SV30 ACETONE	UG/KG	3100	V	39	U	1700	U			
SV31 CARBON DISULFIDE	UG/KG	3100	U	13	U	1700	U	:	:	
SV32 2-BUTANONE	UG/KG	3100	U	13	U	1700	U			
SV33 VINYL ACETATE	UG/KG	NA	0	NA .	0	NA .	0		:	•
SV34 2-HEXANONE	: UG/KG	3100	U	13	U	1700	υ		;	
SV35 4-METHYL-2-PENTANONE	: UG/KG:	3100	บ	13	U	1700	U	:		
SMBG STYRÉHE	: UG/KG:	3100	Ū	: 14	J	1700	ij			
SV3/ AYLENES, TOTAL	: UG/KG:	25000		13	U	1700	IJ,		,	
2701 SAMPLE NUMBER	NA :	011		:012		013	:	101	102	
2202 ACTIVITY CODE	: NA	DSX44		DSX44		DSX44	·····- ·- ·	DSX44	D2X41	

·	COMPOUND	UNITS	103	104	`.	105		106		107	
SMO1 SILVER	BY ICAP	MG/KG: 2.4	U	2.3	Ü	2.3	U	2.4	U	2.2	U
SMO2 ALUMINUM	BY ICAP	MG/KG: 4200		5700		5200		7700		5000	
SMO3 ARSENIC	BY ICAP	MG/KG:5.3		6.0		7.9		5.3		8.1	
SMO4' BARTUM	BY ICAP	MG/KG: 71		110		190		190		86	
SMO5 BERYLLIUM	BY ICAP	MG/KG: 1.2	U	1.1	U	1.1	U	1.2	U	1.1	U
SMO6 CADMIUM	BY ICAP	MG/KG: 1.2	U	1.1	U	1.1	U	1.2	U	1.1	U
SMO7 COBALT	BY ICAP	:MG/KG:12	Ü	11	U	11	U	12	U	11	U
SMO8 CHROMIUM	BY ICAP	MG/KG:7.0		15		11		11		8.2	
SMO9 COPPER	BY ICAP	:MG/KG:13		37		38		19		: 19	 :
SM10 IRON	By ICAP	MG/KG:8300		22000		18000		14000		10000	
SM11 MANGANESE	BY ICAP	:MG/KG:300		380		240		370		230	
SM12 MOLYBDENUM	BY ICAP	:MG/KG:NA	0	NA	0	NA	0	NA	0	NA	0
SM13 NICKEL	BY ICAP	:MG/KG:9.8		13		12		15		9.2	
SM14 LEAD	BY ICAP ;	:MG/KG:66	J	59	J	81	J	59	J	200	J
SM15 ANTIMONY	BY TCAP	MG/KG:14	U	14	U	14	U	14	U	13	Ų
SM16 SELENIUM	BY ICAP	:MG/KG:1.2	IJ	1.1	U	1.1	U	1.3		1 1	U :
SM17 TITANIUM	BY ICAP	MG/KG:NA	0	NA	0	NA	0	NA	0	NA	1 0
SM18 THALLIUM	BY ICAP	MG/KG: 2.4	U	2.3	Ü	2.3	U	2.4	U	2.2	U :
SM19 VANADIUM	BY TCAP	MG/KG 12		20		17	:	20		12	:
5M20 ZTNI	BY TUAP	MG7FG 53	J	110	j	170	J	71	.1	77	
SM21 CALCIUM	BY ICAP	Mit/FG:58000	:	59000	:	26000	:	29000		79000	10
SM22 MAGNESTUM	B's TUAP	MG/FG: 7000		9600	:	2300	:	5000		7600	
SM23 SODIUM	BY ICAP	MG/KG: 1200	U :	1100	U :	1100	U	12000		1100	
SM24 POTASSIUM	B7 ICAP	MG/NG: 1200	U	1100	U	1100	U	1500	: •	1100	u .
SSO1 PHENOI		UG/KG: 10000	U	11000	IJ	12000	U :	360	U :	11000	u
SSO2 CARBAZOLE		UG/KG: 10000	U	11000	U	12000	U	380	U	11000	U
			•				•				

COMPOUND	UNITS	103		104		105	•	10	06	107	, .
SSO3 BIS(2-CHLOROETHYL) ETHER	UG/KG:	10000	U	11000	 U	12000	-	380	U	. 11000	 U
SSO4 2-CHLOROPHENOL	UG/KG	10000	U	11000	U	12000	บ	380	U	11000	U
SSOS 1.3-DICHLOROBENZENE	UG/KG:	10000	U	11000	Ü	12000	U	380	U	11000	U
SSO6 1.4-DICHLOROBENZENE	UG/KG:	10000	U	11000	Ū	12000	U	380	U	11000	Ü
SSO7 BENZYL ALCOHOL	UG/KG:	NA	0	NA	ō	NA	0	NA	0	NA	0
SSOB 1.2-DICHLOROBENZENE	UG/KG:	10000	V	11000	U	12000	U	380	U	11000	U
SSO9 2-METHYLPHENOL (O-CRESOL)	UG/KG:	10000	U	11000	U	12000	U	380	U	11000	U
SS10 BIS(2-CHLOROISOPROPYL) ETHER .	UG/KG:	10000	U	11000	U	12000	U	380	U	11000	U
SS11 4-METHYLPHENOL (P-CRESOL)	UG/KG:	10000	, U	11000	Ü	12000	U	380	 U	11000	U
SS12 N-NITROSO-DIPROPYLAMINE	UG/KG:	10000	U	11000	U	12000	U	380	u	11000	11
SS13 HEXACHLOROETHANE	UG/KG:	0000	U	11000	U	12000	U	380	U	11000	
SS14 NITROBENZENE	UG/KG:	0000	U	11000	บ	12000	U	380	U	11000	U
SS15 ISOPHORONE	UG/KG: 1	0000	U	11000	U	12000	U	380	U	11000	U
SS16 2-NITROPHENOL	UG/KG: 1	0000	U	11000	U	12000	U	380	U	11000	Ü
SS17 2.4-DIMETHYLPHENOL	UG/KG: 1	0000	U	11000	U	12000	U	380	U	11000	. U
SS18 BENZOIC ACID	UG/KG:N	IA	0	NA	0	NA	0	NA	0	NA	0
SS19 BIS(2-CHLOROETHYOXY) METHANE	UG/KG: 1	0000	U	11000	U	12000	U	380	U	11000	, n
SS20 2,4-DICHLOROPHENOL	UG/KG: 1	0000	U	11000	Ü	12000	U	380	U :	11000	U
SS21 1.2.4-TRICHLOROBENZENE	:UG/KG:1	0000	U	11000	U	12000	U	380	11	11000	U.
51-22 NAPH THAT ENE	UG/KG: 1	0000	U	11000	U :	12000	· · · · · · · · · · · · · · · · · · ·	380	1)	P0000	i
5523 4-0 HEORDANILINE	UG/KG: 1	0000	U	11000	U :	12000	U	380	11	14600	11
5521 HEXACHLOROBUTADIENE	UG/KG: 1	0000	U :	11000	U :	12000	Ū.	380	Ü	11000	Ü :
5S25 4- CHLORO-3-METHYLPHENOL	UG/KG: 1	0000	U :	11000	U :	12000	U :	380	11	11000	u u
SS26 2 METHYLNAPHTHALENE	UG/KG: 1	0000	U :	11000	U :	12000	U .	380	บ	1 3000	. <u></u>
5527 HEXACHLOROCYCLOFENTADIENE	UG/KG: 10	0000		11000	U	12000	U	380	U ,	11000	11
SS28 2.4.6-TRTCHLOROPHENOL	UG/KG: 10	0000	U	11000	Ü	12000	Ü	380	<u>U</u>	11000	u :

COMPOUND	UNITS	1	03	104		105		106		107	7
SS29 2.4.5-TRICHLOROPHENOL	UG/KG	54000	บ	58000	U	60000		2000	U	57000	U
SS30 2-CHLORONAPHTHALENE	:UG/KG	10000	U	11000	U	12000	U	: 380	U	11000	Ü
SS31 2-NITROANILINE	UG/KG	54000	U	58000	U	60000	U	2000	U	57000	U
SS32 DIMETHYLPHTHALATE	UG/KG	10000	U	11000	U	12000	U	380	U	11000	U
SS33 ACENAPHTHYLENE	UG/KG	10000	U	11000	U	12000	U	380	v	11000	U
SS34 3-NITROANILINE	UG/KG:	54000	U	58000	U	60000	U	2000	U	57000	U
SS35 ACENAPHTHENE	UG/KG:	10000	U	11000	U	12000	U	: 380	U	11000	U
SS36 2.4-DINITROPHENOL	UG/KG	54000	v	58000	U	60000	U	2000	U	57000	บ
SS37 4-NITROPHENOL	UG/KG	54000	U	58000	U	60000	U	2000	U	57000	U
SS38 DIBENZOFURAN	: UG/KG:	10000	U	11000	U	12000	U	: 380	ij	11000	บ
SS39 2.4-DINITROTOLUENE	UG/KG:	10000	U	11000	U	12000	U	380	U	11000	Ü
SS40 2,6-DINITROTOLUENE	UG/KG:	10000	U	11000	U	12000	V	380	U	11000	U
SS41 DIETHYLPHTHALATE	UG/KG	10000	U	11000	U	12000	U	380	U	11000	U
SS42 4-CHLOROPHENYL PHENYL ETHER	UG/KG	10000	U	11000	U	12000	· U	380	U	11000	U
SS43 FLUORENE	UG/KG:	10000	U	11000	U	12000		380	U	11000	ń.
SS44 4-NITROANILINE	:UG/KG:	54000	U	58000	U	60000	U	2000	U :	57000	Ū:
SS45 4.6-DINITRO-2-METHYLPHENOL	UG/KG:	54000	U	58000	U	60000	U	2000	U	57000	-U
SS46 N-NITROSODIPHENYLAMINE	UG/KG:	10000	U	11000	U	12000	U	380	U	11000	U :
SS47 4-BROMOPHENYL PHENYL ETHER	UG/KG:	10000	U	11000	<u> </u>	12000	U	380	U	11000	<u>U</u>
State HEXACHLOROBENZEM	. 06/66	10000	ti :	11000	U	12000	·	380	()	11000	· · · · · · · · · · · · · · · · · · ·
S549 PENTACHI OROPHENOL	:UG/KG:	0000	U	58000	Ü	60000	0	S000	Ü	57000	U
SSSO PHEMANTHRENE	UG/FG:	10000	U	20000	:	55000		1500		11000	U
SSS1 ANTHRACENE	:UG/KG:	0000	Ü :	11000	U :	15000	:	380	U	11000	Ü
SS52 DI H-BUTYL PHTHALATE	UG/KG:	0000	U	11000	U	12000	U :	380	U	11000	ii .
SS53 FLUORANTHENE	UG/KG: 1	10000	Ū	4000 .	:	38000	:	2000	 	11000	U :
SS54 PYRENE	UG/KG	3000	1	50000		83000	us ids	2200 Jil	311	11000	U

COMPOUND UNITS 103 104 106 107 105 SS55 BUTYL BENZYL PHTHALATE UG/KG: 10000 U :11000 U : 12000 U : 386 U : 11000 SS56 3,3'-DICHLOROBENZIDINE :UG/KG: 10000 U:11000 U:12000 U :380 U:11000 U U \$21000 SS57 BENZO(A)ANTHRACENE : UG/KG: 10000 :28000 :1300 :11000 U SS58 BIS(2-ETHYLHEXYL)PHTHALATE : UG/KG: 10000 U :11000 U:12000 U :380 U:11000 U : S\$59 CHRYSENE U 329000 UG/KG: 10000 : 29000 : 1200 :11000 SS60 DI-N-OCTYL PHTHALATE UG/KG: 10000 U:11000 U :12000 U :380 U:11000 SS61 BENZO(B)FLUORANTHENE UG/KG: 10000 U :17000 U : U :11000 : 1000 :11000 SS62 BENZO(K)FLUORANTHENE : UG/KG: 10000 U :11000 U \$12000 . :890 :11000 SS63 BENZO(A)PYRENE UG/KG: 10000 U :**₹6000** 22000 :1100 :11000 U SS64 INDENO(1,2.3-CD)PYRENE UG/KG: 10000 U:11000 U :12000 U :660 :11000 SS65 DIBENZO(A.H)ANTHRACENE : UG/KG: 10000 U:11000 U:12000 U :380 U:11000 SS66 BENZO(G.H.I)PERYLENE UG/KG: 10000 U:11000 U 2000 :710 £ 11000 5.9 U \$14 STO9 CYANIDE MG/KG:94 : 220 : 190 2201 SAMPLE NUMBER :106 : NA : 103 : 104 : 105 : 107 2202 ACTIVITY CODE **DSX44** :DSX44 NA DSX44 : DSX44 : DSX44

٠		COMPOUND	UNITS	108	109		110		111		112	
SMO1 SI	ILVER E	Y ICAP	MG/KG	2.6	2.4	U	2.5	U	2.5	ບ	2.7	· . U
SMO2 AL	UMINUM 8	Y ICAP	MG/KG	5200	6600		11000		3800		7600	
SMO3 AR	RSENIC B	Y ICAP	MG/KG	14	11		10		7.8		33	
SMO4 BA	ARIUM B	Y ICAP	MG/KG	160	150		220		96		170	
SMO5 BE	RYLLIUM B	Y ICAP	MG/KG	1.3 L	1.2	U	1.2	U	1.2	U	1.3	U
SMO6 CA	OMIUM B	Y ICAP	MG/KG:	8.2	1.2	U	1.2	U	1.2	U	1.3	U
SMO7 CO	BALT B	Y ICAP	MG/KG	13 U	12	U	12	U	12	U	13	U
SMO8 CH	ROMIUM B	Y ICAP	MG/KG	20	15		23		15		12	
SMO9 CO	PPER B	Y ICAP	MG/KG	270	93		56		72		43	
SM10 IRO	ON B	Y [I.AP	MG/KG	21000	28000		29000		33000		24000	
SM11 MAN	NGANESE B	Y ICAP	MG/KG:	180	390		470	J	320	J	870	J
SM12 MOL	LYBDENUM B	/ ICAP	MG/KG:	NA O	: NA	0	NA	0	: NA	0	: NA	0
SM13 NIC	CKEL 8'	/ ICAP	MG/KG:	14	:20		23		27		23	
SM14 LEA	AD B	/ ICAP	MG/KG:	270 J	230	J	230	J	150	J	300	J
SM15 ANT	TIMONY BY	/ ICAP	MG/KG	15 U	:14	U	15	U	: 15	U	16	U :
SM16 SEL	LENIUM BY	/ ICAP	MG/KG:	2.3	1.6		3.0	J	1.9		1.5	J
SM17 TIT	TANIUM BY	' ICAP	MG/KG:	NA O	NA	0	NA	0	NA	0	NA	C
SM18 THA	ALTUM BY	ICAP	MG/KG:	2.6 U	2.4	U	2.5	U	2.5	U	2.7	Ú
SM19 VAN	VADIUM BY	ICAP	MG/KG:2	25	23		32		17		23	
5M20 Z1N	lı RA	II AF	MG/KG:2	2500 J	550	J	360		250		530	
M21 (4)	C1000 B7	II AP	MG/FG: 2	2700	33000		28000		3600		37000	• • •
SM22 MAG	JIESTUM BY	TUAP	MG/KG: 1	300 U	2300	:	3300		1200	U	6400	-
SM23 SOD	DIUM BY	II AP	MG/KG: 1	300 U	1200	U	1200	U	1200	Ü	1300	ij
SM2-I POT	ASSIUM BY	ICAP	MG/KG 1	300 U	1200	U	1200		1200	U	1300	
SSOT PHE	NOL		UG/KG:3	390 U	420	U	390	U	390	Ü	410	Ü
5502 CAR	RBAZOLE		UG/KG 4	120 J	500	:	390	U	390	Ü :	410	U :

COMPOUND	UNITS	108	109		110	111	112
SSO3 BIS(2-CHLOROETHYL) ETHER	UG/KG	390 U	420	V	390 U	:390 U	.410 U
SSO4 2-CHLOROPHENOL	UG/KG	390 U	420	U	390 U	:390 U	:410 U
SSOS 1.3-DICHLOROBENZENE	UG/KG	390 U	420	U	390 U	:390 U	410 U
SSO6 1.4-DICHLOROBENZENE	UG/KG	390 U	420	U	390 U	:390 U	410 U
SSO7 BENZYL ALCOHOL	UG/KG	NA O	NA	0	NA O	NA O	NA O
SSO8 1.2-DICHLOROBENZENE	UG/KG:	390 U	420	U	390 U	390 บ	410 U
SSO9 2-METHYLPHENOL (O-CRESOL)	UG/KG:	390 U	420	U	390 U	390 U	410 U
SS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/KG	390 U	420	U	390 U	390 U	410 U :
SS11 4-METHYLPHENOL (P-CRESOL)	UG/KG	390 U	420	U	390 U	390 U	410 U
SS12 N-NITROSO-DIPROPYLAMINE	. UG/KG:	390 U	420	U	390 U	390 U	410 II
SS13 HEXACHLOROETHANE	UG/KG:	390 U	: 420	U :	390 U	390 U	410 U
SS14 NITROBENZENE	UG/KG	390 U	420	U	390 U	390 U	410 U
SS15 ISOPHORONE	UG/KG	390 U	420	Ū	390 U	390 U	410 U
SS16 2-NITROPHENOL	UG/KG:	390 U	420	Ū	390 U	390 U	410 U
SS17 2.4-DIMETHYLPHENOL	UG/KG	390 U	420	Ū	390 U	390 (/	410 4
SS18 BENZOIC ACID	UG/KG:	NA O	:NA (0 :1	NA O	NA O	NA O
SS19 BIS(2-CHLOROETHYOXY) METHANE	UG/KG:3	390 U	420	Ū	390 U	390 U	410 U
SS20 2.4 DICHLOROPHENOL	UG/KG:3	390 U	420 l	U	390 U	390 . U	410 U
SS21 1.2.4-TRICHLOROBENZENE	UG/KG:3	390 U	: 420 L	U ::	390 U	390 ()	410 U
5-22 NAPHTHALENE	804500	100	510	:	390 U	390 #	410
\$523 4-OROROATH THE	0G/KG :	390 1)	: 420	v : 3	390 0	390 0	410 0 1
SS24 HEXACHI OROBUTAD LEHE	UGZEG: 3	390 ()	: 120	u : 3	390 "	390 11	410
SS25 4-CHLORO 3-METHYLPHCHOL	.UG/kG.3	390 U	420 U	U : 3	390 U	390 U	410 U
SS26 2 METHYLNAPHTHALENE	UG/KG	2500	: 420 U	נו נו	390 U	390 U	410 U
SS27 HEXACHLOROCYCLOPENTADIENE	UG/KG: 3	90 U	:420 U	3 3	390 U	390 U	410 (1
5528 2.4.6-TRTCHLOROPHENOL	UG/KG:3	190 U	420 U	J : 3	390 U	390 0	410 0
	: :-		,	:-	:		

COMPOUND	UNITS	108	109		110	111	112	
SS29 2.4.5-TRICHLOROPHENOL	: UG/KG: 940		: 1000	u	950 U	:940 U	990	U
SS30 2-CHLORONAPHTHALENE	UG/KG: 390	U	420	บ	390 U	390 U	410	U
SS31 2-NITROANILINE	UG/KG: 940	U	1000	U	950 U	.940 U	990	U
SS32 DIMETHYLPHTHALATE	UG/KG: 390	U	420	Ū	390 U	390 U	410	U
SS33 ACENAPHTHYLENE	UG/KG 2400		460		390 U	:390 U	410	U
SS34 3-NITROANILINE	UG/KG 940	U	1000	U	950 U	940 U	990	U
SS35 ACENAPHTHENE	UG/KG 690		420	U	390 U	390 U	410	U
SS36 2.4-DINITROPHENOL	UG/KG 940	U	1000	U	950 U	940 U	.990	U
SS37 4-NITROPHENOL	UG/KG:940	, U	1000	U	950 U	940 U	990	U
SS38 DIBENZOFURAN	UG/KG: 1100		420	U	390 U	390 U	410	li .
SS39 2.4-DINITROTOLUENE	UG/KG: 390	U	420	U	390 U	390 U	410	U :
SS40 2.6-DINITROTOLUENE	UG/KG: 390	Ü	420	U	390 U	390 U	410	Ü
SS41 DIETHYLPHTHALATE	UG/KG: 390	U	420	U	390 U	390 U	410	U
SS42 4-CHLOROPHENYL PHENYL ETHER 1.	UG/KG: 390	U	420	U	390 U	390 U	410	Ü
SS43 FLUORENE	UG/KG 3100	ا المان الرابع المان المان المان المان المان المان المان المان المان المان المان المان المان المان المان المان ا	460	J	390 U	390 11	410	0
SS44 4-NITROANILINE	UG/KG:940	V	1000	U	950 V	940 U	:990	U
SS45 4.6-DINITRO-2-METHYLPHENOL	UG/KG: 940	U	1000	U	950 U	940 U	990	U
SS46 N-NITROSODIPHENYLAMINE	UG/KG 390	U	420	U	390 U	390 U	410	U
SS47 4-BROMOPHENYL PHENYL ETHER	UG/KG 390	U	420	U :	390 U	390 U	410	U
5546 HEXACHI OROBENZENE	UG/KG: 390	IJ	420	:	390 "	390 0	410	ıı.
SS49 PENTACHLOROPHENOL	.UG/KG: 940	U	1000	U	950 U	940 0	ากษ	u
SSSO PHENANTHRENE	UG/KG 43000	J	2200	:	1500	850	: 110	
SSS1 ANTHRACENE	UG/KG: 4700	J	780		390 U	390 11	410	U U
SSS2 DI N-BUTYL PHTHALATE	UG/KG: 390): : U	420	U :	390 U	390 ()	.410	Ü.
5353 FLUORANTHENE	UG/KG: 40000	J	2800	::	2200	770	1300	
SS54 PYRENE	UG/KG: 58000	J	700		2300	81Q 📆	1400	1

COMPOUND	UNITS	10	8	109	r	110)	11	١	112	2
SS55 BUTYL BENZYL PHTHALATE	: UG/KG	: 390	U	420	U	: 390	U	390	U	410	U
SS56 3.3'-DICHLOROBENZIDINE	: UG/KG	390	U	420	U	390	U	390	U	410	U
SS57 BENZO(A)ANTHRACENE	UG/KG	34000	J	500		1500		480		790	
SSS8 BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG	390	U	420	U	390	U	390	U	410	U
SS59 CHRYSENE	UG/KG	34000		4300		1600		550	:	850	
SS60 DI-N-OCTYL PHTHALATE	UG/KG	390	U	420	U	390	U	390	Ū	410	U
SS61 BENZO(B)FLUORANTHENE	UG/KG	16000	J	4900		1100		440		610	
SS62 BENZO(K)FLUORANTHENE	UG/KG	390	υ	3400		1200		450		680	
SS63 BENZO(A)PYRENE	UG/KG	20000		1200		1300		440		700	
SS64 INDENO(1,2,3-CD)PYRENE	: UG/KG	12000	J	2700		760		390	U	410	U
SS65 DIBENZO(A.H)ANTHRACENE	UG/KG	390	U	420	U	390	U	390	U	410	U
SS66 BENZO(G,H,I)PERYLENE	UG/KG	14000	J	2600		760	1 2 11	390	U	410	U
STO9 CYANIDE	MG/KG	98	1.50	35		6.2	U	6.2	U	6.7	U
ZZO1 SAMPLE NUMBER	. NA	108		109		110		111		112	
ZZOZ ACTIVITY CODE	NA	DSX44		DSX44		DSX44		DSX44		D5X44	•

WFO1 WATER TEMP 'C 11.0 7.0 9.0 9.0 WFO5 PH. FIELD SU 7.58 7.81 6.32 6.21	
WF10 CONDUCTIVITY (FIELD)	 U
WMO1 SILVER	U
WMO2 ALUMINUM BY ICAP UG/L 45000 J :15000 J :200 U :200 U :200 WMO3 ARSENIC BY ICAP UG/L 40 670 15 J :10 U :10 WMO4 BARIUM BY ICAP UG/L 2100 890 1000 630 200 WMO5 BERYLLIUM BY ICAP UG/L 5.0 U :5.0 U :5	U
WMO3 ARSENTC BY ICAP UG/L 40 670 15 J 10 U 10 WMO4 BARIUM BY ICAP UG/L 2100 890 1000 630 200 WMO5 BERYLLIUM BY ICAP UG/L 5.0 U	
WMO4 BARIUM BY ICAP UG/L 2100 890 1000 630 200 WMO5 BERYLLIUM BY ICAP UG/L 5.0 U	U
WMO5 BERYLLIUM BY ICAP UG/L 5.0 U >U</td>	U
WMO6 CADMIUM BY ICAP UG/L 5.0 U 5.0 U 19 21 5.0 WMO7 COBALT BY ICAP UG/L 50 U	U
WMO7 COBALT BY TCAP UG/L 50 U 10 U <	U
WMO8 CHROMIUM BY ICAP UG/L 68 29 10 U 10 U 10 WM09 COPPER BY ICAP UG/L 84 J 91 J 37 U 27 U 200 WM10 IRON BY ICAP UG/L 110000 53000 30000 15000 U 320 WM11 MANGANESE BY ICAP UG/L 5600 850 1300 630 32 WM12 MOLYBDENUM BY ICAP UG/L NA O	U
WMO9 COPPER BY ICAP UG/L 84 J 91 J 37 U 27 U 200 WM10 IRON BY ICAP UG/L 110000 53000 30000 15000 U 320 WM11 MANGANESE BY ICAP UG/L 5600 850 1300 630 32 WM12 MOLYBDENUM BY ICAP UG/L NA O NA O NA O NA O NA O N/A WM13 NICKEL BY ICAP UG/L 95 40 U 60	U
WM10 IRON BY ICAP UG/L 110000 53000 30000 15000 U 320 WM11 MANGANESE BY ICAP UG/L 5600 850 1300 630 32 WM12 MOLYBDENUM BY ICAP UG/L NA O NA O NA O NA O NA O N/A WM13 NICKEL BY ICAP UG/L 95 40 U 40	IJ
WM11 MANGANESE BY ICAP UG/L 5600 850 1300 630 32 WM12 MOLYBDENUM BY ICAP UG/L NA O ></td>	
WM12 MOLYBDENUM BY ICAP UG/L NA O	U
WM13 NICKEL BY ICAP UG/L 95 40 U :40 U :40 U :40 WM14 LEAD BY ICAP UG/L :58 1200 J :330 350 3 6 WM15 ANTIMONY BY ICAP UG/L :60 U :60 U :60 U :60 U :60	U
WM14 LEAD BY ICAP UG/L 58 1200 J 330 350 3 6 WM15 ANTIMONY BY ICAP UG/L 60 U 60 U 60 U 60 U 60	n,
WM15 ANTIMONY BY ICAP UG/L :60 U :60 U :60 U :60	U
	O
WM16 SELENTUM BY ICAP UG/L :25 U :25 U :NA I :NA I :N/A	V
	I
WHITE TELEMENT OF THE TELEMENT	Ú
WHITE THOLLIUM BY LEAP UG/L: 10 U:50 U:NA I:NA I N/A	1
WELLS VARIABLEM BY TEAP UG/L:77 :50 U:50 U:50 U:50	U
WM20 21NC BY 10AP .UG/L :4200 :1100 :40000 :35000 -41	
WHI21 CALCIUM. TOTAL BY ICAP MG/L .520 320 210 U 230 U 50	
WM22 MAGHESTUM. TOTAL BY ICAP MG/L:66 :44 :37 U:34 U:19	
WM23 SODIUM TOTAL BY ICAP MG/L:110:150:99:64 U:19	

CON	POUND	UNITS	201	202		203	204	205F
WM24 POTASSIUM, TOTAL BY	' ICAP	MG/L	53	. 46		: 33	26	:5.0
WM35 SILVER, DISSOLVED	BY ICAP	UG/L	10 U	10	U	10 U	10 U	:
WM36 ALUMINUM.DISSOLVED	BY ICAP	UG/L	200 U	200	υ	:200 U	200 U	:
WM37 ARSENIC.DISSOLVED	BY ICAP	UG/L	13	80	J	14 U	10 U	:
WM38 BARIUM.DISSOLVED	BY ICAP	UG/L	490	260	``	650	490	
WM39 BERYLLIUM.DISSOLVED	BY ICAP	UG/L	5.0 U	5.0	U	5.0 Ū	5.0 U	:
WM40 CADMIUM.DISSOLVED	BY ICAP	UG/L	5.0 U	5.0	U	5.0 U	5.0 U	: :
WM41 COBALT.DISSOLVED	BY ICAP	UG/L	50 U	50	U	.50 U	50 U	:
WM42 CHROMIUM, DISSOLVED	BY ICAP	UG/L	10 U	10	U	:10 U	10 U	:
WM43 COPPER DISSOLVED	BY ICAP	UG/L	25 U	25	U	25 U	25 U	
M44 IRON.DISSOLVED	BY ICAP	UG/L	1500 J	14000	J	20000	11000	:
M45 MANGANESE DISSOLVED	BY ICAP	UG/L	1500	540		1100	600	
M46 MOLYBDENUM.DISSOLVE	D BY ICAP	UG/L	NA O	NA	0	NA O	NA O	
M47 NICKEL DISSOLVED	BY ICAP	:UG/L	40 U	40	U	34 U	40 U	
M48 LEAD DISSOLVED	BY ICAP	UG/L	5 0 J	3.0	U	3.8 U	5 3 U	
M49 ANTIMONY, DISSOLVED	BY ICAP	:UG/L	60 U	60	U	60 U	60 U	
M50 SELENIUM,DISSOLVED	BY ICAP	UG/L	5.0 U	5.0	U	5.0 U	20 U	
M51 TITANIUM, DISSOLVED	BY ICAP	UG/L	NA O	:NA	0	NA O	NA O	
M52 THALLIUM, DISSOLVED	BY ICAP	UG/L	10 U	50	U	10 ປ	10 U	
M53 VANADIUM DISSOLVED	BY ICAP	UG/I	50 U	50	U	50 11	50 U	
Mod Zinc Dissolven	BY TEAP	HG/L	44	:23	: ::	690	3400	
M55 CALCIUM, DISSOLVED	BY ICAP	MG/L	170	270		. 220	250	
MS6 MAGNESTUM.DISSOLVED	BY ICAP	MG/L .	41	: 42	-·- :	39 U :	35	
MIST SODIUM, DISSULVILD	BY ICAP	MG/L :	100	. 150	: :	100	66	
MS8 POTASSTUM.DISSOLVED	BY ICAP	MG/L	33	: 44	: - - : :	NA O	26	
SO1 PHENOL	*********	UG/L :	10 U	10	U :	10 0	10 ປ	10 1

COMPOUND	UNITS	.	201		202	203	204		205F
WSO3 BIS(2-CHLOROETHYL) ETHER	:UG/L	10	u	10	U	: 10 U	: 10	10	 U
WSO4 2-CHLOROPHENOL	UG/L	10	U	10	U	10 U	10 l	10	U
WSO5 1.3-DICHLOROBENZENE	:UG/L	10	U	10	U	10 U	10	10	U
WSO6 1.4-DICHLOROBENZENE	UG/L	10	U	10	U	10 U	10	10	U
WSO7 BENZYL ALCOHOL	UG/L	10	U	10	U	10 U	10	10	U
WSOB 1.2-DICHLOROBENZENE	UG/L	10	U	10	U	10 U	10 0	10	V
WSO9 2-METHYLPHENOL (O-CRESOL)	UG/L	10	U	10	U	10 U	10 U	10	U :
WS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/L	10	U	10	U	10 U	10 ປ	10	U
WS11 4-METHYLPHENOL (P-CRESOL)	UG/L	10	U	10	_ U	10 U	10 U	10	U :
WS12 N-NITROSO-DIPROPYLAMINE	UG/L	10	U	10	U	: 10 U	10 U	10	1)
WS13 HEXACHLOROETHANE	UG/L	10	U	: 10	U	10 U	:10 U	10	บ :
WS14 NITROBENZENE	UG/L	10	U	: 10	U	10 U	10 U	10	U :
WS15 ISOPHORONE	UG/L	10	U	10	U	10 U	10 U	10	U
WS16 2-NITROPHENOL	UG/L	10	V	10	U	10 U	:10 U	10	U
WS17 2.4-DIMETHYLPHENOL	UG/L	10	U	10	U	10 U	: 10 U	10	: U
WS18 BENZOIC ACID	UG/L	50	U	50	U	50 U	:50 U	50	U :
WS19 BIS(2-CHLOROETHYOXY) METHANE	UG/L	10	U	10	U	10 U	10 U	10	-1J
WS20 2,4-DICHLOROPHENOL	UG/L	10	U	10	U	10 U	: 10 U	. 10	U .
WS21 1.2.4-TRICHLOROBENZENE	UG/L	10	U	10	U	10 U	: 10	10	U
WSZZ NAPITHALENE	UG/L:	10	U	10	U :	10 U	10 0	10	U
W523 4-4 HLURQANTI INE	UG/L :	10	U	10	U	10 U	10 0	10	u
WS24 HEXACHLOROBUTADIENE	UG/L :	10	U	10	U	10 U	10 11	10	U
W525 4 CHLORO 3-METHYLPHENOL :	UG/L:	10	U	10	U :	10 U	10 U	10	IJ
WS2G 2 METHYLNAPHTHALENE	UG/L :	10	U :	10	U :	10 U	10 U	10	Ü
WS27 HEXACHLOROCYCLOPENTADIENE	UG/L:	10	v	10	V	10 U	10 U	10	Ū
WS28 2,4.6-TRTCHLOROPHENOL	UG/L	10	U	10	U	10 U	10 ป	10	Ü

	COMPOUND	UNITS	•	201		202		203		204		205F
WS29 2.4.	.5-TRICHLOROPHENOL	UG/L	: 25	U	: 25	U	: 25	U	25	U	25	U
WS30 2-CH	ILORONAPHTHAL ENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS31 2-NI	TROANILINE (ORTHO NITROANILINE)	UG/L	25	U	: 25	U	25	V	: 25	U	: 25	U
WS32 DIME	THYLPHTHALATE	UG/L	10	U	:10	U	10	U	10	U	10	U
WS33 ACEN	IAPHTHYLENE	UG/L	10	U	10	U	10	U	10	U	10	V
WS34 3-NI	TROANILINE	UG/L	25	V	:25	U	25	U	25	U	25	U
WS35 ACEN	IAPHTHENE	UG/L	10	V	10	U	:10	U	65		10	U
WS36 2.4-	DINITROPHENOL	UG/L	25	U	25	U	25	U	25	V	25	U
WS37 4-NI	TROPHENOL	UG/L	25	U	25	U	25	U	25	U	25	U
WS38 DIBE	NZUFURAN	UG/L	10	U	10	U	10	U	10	U	10	U
WS39 2.4-0	DINITROTOLUENE	UG/L	10	U	10	U	10	U	10	U	10	. U
WS40 2.6-	DINITROTOLUENE	UG/L	10	U	:10	U	10	U	10	U	10	U
WS41 DIETI	HYLPHTHALATE	UG/L	10	U	10	U	10	U	10	U	10	U
WS42 4-CH	LOROPHENYL PHENYL ETHER	UG/L	10	U	10	U	10	U	10	U	10	U
WS413 FLUOR	RENE	UG/L	10	U	10	U	10	U	25		10	ij.
WS44 4-NI1	TROANILINE	UG/L	25	U	25	U	25	Ü	25	U	25	U
WS45 4.6-0	DINITRO-2-METHYLPHENOL	UG/L	25	U	25	U	25	U	25	U	25	1)
WS46 N-N11	TROSODIPHENYLAMINE	UG/L	10	V	10	U	10	U :	10	U	10	U
WS47 4-BRC	OMOPHENYL PHENYL ETHER	UG/L	10	U	10	U	10	U :	10	U	10	U
World 111 Add	मा लेखित छ हो ।।	11671	10	Ų	10	U	10	U :	10		: 10	U
w as PENDA	No fact officially from	UG/L	25	U	25	U	25	ŧi :	25	U	25	Ü
WSCO PHENA	MITHRENE	UG/L	10	U	10	U	10	U	16	· · · · · · · · · · · · · · · · · · ·	10	U
WS51 ANTHR	RACENE	UG/L	10	U	10	U	10	U	10	IJ	10	(1)
WSC2 DI H-	BUTYL PHILIALATE	UG/L .	10	U	10	U	10	U	10	U	10	U .
WS53 FLUOR	RANTHENE	UG/L	10	U	10	U	10	U :	10	Ü	10	II ·
WS54 PYREN	IF	UG/L	10	U	10	U	10	U :	10	U	10	U

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-DSX44

COMPOUND	UNITS	•	201		202	203	3	2	204	20	5f
WS55 BUTYL BENZYL PHTHALATE	UG/L	:10	U	:10	v	:10	U	10	U	. 10	U
WS56 3.3'-DICHLOROBENZIDINE	UG/L	10	U	10	V	10	U	10	U	10	U
WS57 BENZO(A)ANTHRACENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/L	10	U	:10	U	10	U	10	U	10	U
WS59 CHRYSENE	UG/L	10	U U	10	V	10	U	10	U	10	U
WS60 DI-N-OCTYL PHTHALATE	UG/L	10	U	10	U	10	U	10	U	10	U
WS61 BENZO(B)FLUORANTHENE	UG/L	10	U	10	V	10	U	10	U	10	U
WS62 BENZO(K)FLUORANTHENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS63 BENZO(A)PYRENE	UG/L	10	Ų	10	U	10	U	10	U	10	U
WS64 INDENO(1.2.3-1D)PYRENE	UG/L	10	V	10	υ	10	U	10	U	10	υ :
WS65 DIBENZO(A,H)ANTHRACENE	UG/L	10	U	10	U	10	U	10	(I	10	U
WS66 BENZO(G.H.I)PERYLENE	UG/L	10	U	10	U	10	U	10	U	10	U :
WS67 CARBAZOLE	UG/L	10	U	10	U	10	U	10	U	10	U
WTO9 CYANIDE, TOTAL	MG/L	0.52	J	0.027	J	0.590	J	1.6	J	0.013	J
WVO3 CHI OROMETHANE	UG/L	10	U	10	U	10	U	10	U.	10	0
WVO4 BROMOME THANE	UG/L	10	U	10	U	10	U	10	U	: 10	U
WVO5 VINYL CHIORIDE	UG/L	10	ป	10	U	10	Ü	10	ป	10	U
WVO6 CHLOROETHANE	UG/L	10	U	10	U	10	U	10	U	10	U .
WVO7 METHYLENE CHLORIDE	UG/L	10	U	10	U	10	U	10	U	10	u
WVO3 1 + D1CHLOROETHEBE	11671	10	U	10	U :	10	IJ	10	IJ	10	U
WVOS 1.1-DTI HI OROETHAN:	10671	10	U	10	υ :	10	υ	10	11	10	11
W/10 1 2 DICHLOROETHERE. TOTAL	:06/L :	10	U :	10	U .	10	Ū	10	Ü	10	II
WV11 CHLOROFORM	UG/L	10	U	10	U	10	U	10	U	17	
WV12 1.2 DICHLORGETHANE	UG/L .	10	U	10	U	10	IJ	10	U	10	U ,
WV13 1.1 1-TRICHLOROETHANE	UG/L	10	U	10	U	10	Ü	10	U .	10	Ü
WV14 CARBON TETRACHLORIDE	UG/L	10	U	10	U :	10	U	10	U :	10	U i

COMPOUND	UNIT	5	201		202		203	204		205	F
WV15 BROMODICHLOROME THANE	UG/L	:10	U	10	U	10	U	10	U	.10	U .
WV16 1,2-DICHLOROPROPANE	UG/L	10	U	10	V	10	V	10	U	10	U
WV17 BENZENE	UG/L	10	V	10	U	10	U	93		10	U
WV19 TRICHLOROETHENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV20 CIS-1.3-DICHLOROPROPENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV21 DIBROMOCHLOROMETHANE	UG/L	10	U	10	บ	10	บ	10	U	10	บ
WV22 1,1,2-TRICHLOROETHANE	UG/L	10	U	10	U	10	U	10	Ú	10	U
WV24 BROMOFORM	UG/L	10	U	10	U	10	U	10	U	: 10	U
WV25 TETRACHLOROETHENE	UG/L	10	U	10	U	10	U	10	U	10	U .
WV26 TOLUENE	UG/L	10	· U	10	U	10	U	10	V	10	Ū
WV27 1,1.2.2-TETRACHLOROETHANE	UG/L	10	U	10	U	10	U	10	ti	10	U
WV28 CHLOROBENZENE	UG/L	10	U	10	บ	10	V	10	V	10	U :
WV29 ETHYL BENZENE	UG/L	10	υ	10	U	10	υ	10	. U	10	υ :
WV30 ACETONE	UG/L	10	U	10	U	10	U	10	U	10	U :
WV31 CARBON DISULFIDE	:UG/L	10	U	10	U	10	U	10	11	10	u :
WV32 2-BUTANONE	UG/L	10	U	10	U	10	U	10	U	:10	Ü
WV33 VINYL ACETATE	:UG/L	10	U	10	U	10	U	10	U	10	U
WV34 2-HEXANONE	:UG/L	10	U	10	U	10	U	10	U	: 10	U
WV35 4-METHYL-2-PENTANONE	:UG/L	10	V	10	U	10	U	10	U	: 10	U :
WV 16 STYRETH	UG/L	10	U	10	U	10	0	10	li .	10	11
WV3/ XYLENES. TOTAL	UG/L	10	U	10	U :	10	(1 -	10	· u	10	ü
WY40 TRANS- 1.3-DICHLOROPROPENE	UG/L	10	Ü	10	U	10	U	10	Ü	10	U
2201 SAMPLE NUMBER	: NA	201		202	· • • • • • • • • • • • • • • • • • • •	203		204		205	
ZZGZ ALTIVITY CODE	. NA	DSX44		DSX44	 	DSX4	4	D5X44		DSX44	

COMPOUND	UNITS	206	206D	207F	208F	209F
WM24 POTASSIUM, TOTAL BY ICAP	MG/L	12 U	: 12 U	5.0 U	5.0 U	·;
WM35 SILVER, DISSOLVED BY ICAP	UG/L	10 U	10 U	10 U	10 U	:
WM36 ALUMINUM.DISSOLVED BY ICAP	UG/L	200 U	200 U	200 U	200 U	
WM37 ARSENIC.DISSOLVED BY ICAP	UG/L	29 J	:41	10 U	10 U	
WM38 BARIUM, DISSOLVED BY ICAP	:UG/L	400	440	200 U	200 U	
WM39 BERYLLIUM, DISSOLVED BY ICAP	:UG/L	.5 U	5.0 U	:5.0 U	:5.0 U	
WM40 CADMIUM.DISSOLVED BY ICAP	UG/L	.5 U	5.0 U	່5.0 ປ	5.0 U	
WM41 COBALT, DISSOLVED BY ICAP	UG/L	50 U	50 U	:50 U	.50 U	:
WM42 CHROMIUM.DISSOLVED BY ICAP	UG/L	:10 U	10 U	10 U	10 U	
WM43 COPPER.DISSOLVED BY ICAP	UG/L	25 U	:25 U	25 U	. 25 ປ	
WM44 IRON.DISSOLVED BY ICAP	UG/L	24000 U	27000	: 100 U	: 100 0	:
WM45 MANGANESE.DISSOLVED BY ICAP	UG/L	380	400	15 U	15 U	
WM46 MOLYBDENUM.DISSOLVED BY ICAP	:UG/L	NA O	NA O	NA O	N/A O	::
WM47 NICKEL, DISSOLVED BY ICAP	UG/L	40 U	40 U	40 U	40 U	:
WM48 LEAD DISSOLVED BY ICAP	UG/L	4 0 U	4.8 U	3.0 U	3 0 11	:
WM49 ANTIMONY, DISSOLVED BY ICAP	UG/L	60 U	60 U	:60 U	60 U	
WM50 SELENIUM, DISSOLVED BY ICAP	UG/L	NA I	20 U	5.0 U	5.0 U	
WM51 TITANIUM, DISSOLVED BY ICAP	UG/L	NA O	NA O	NA O	N/A O	
WM52 THALLIUM.DISSOLVED BY ICAP	UG/L	NA I	10 ป	. 10 U	10 "	
WMSG VANADIUM DISSOLVED BY ICAP	UG/I	50 U	50 1/	50 U	50 0	
WM54 21NC.DESSOLVED BY ICAP	UG/L	2000	3200 J	20 0	20 0	
WMS5 CALCIUM, DISSOLMED BY ICAP	MG/L	220 U	240	.5.0 U	50 0	
WMSG MAGNESTUM, DISSOLVED BY ICAP	MG/L	51 U	55	:5.0 U	5.0	
WM57 SODIUM.DISSOLVED BY ICAP	MG/L	230	230	5.0 U	5 0 U	
WMG8 POTASSIUM.DISSOLVED BY ICAP	MG/L	12 U	12	5.0 U	5.0 11	
WSO1 PHENOL	UG/L	10 U	10 U	10 U	10 U	

WSO5 1.3-DICHLOROBENZENE :UG/L :10 U :10 WSO6 1.4-DICHLOROBENZENE :UG/L :10 U :10 WSO7 BENZYL ALCOHOL :UG/L :10 U :10 WSO8 1.2-DICHLOROBENZENE :UG/L :10 U :10	U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	
WSO5 1.3-DICHLOROBENZENE :UG/L :10 U :10 WSO6 1.4-DICHLOROBENZENE :UG/L :10 U :10 WSO7 BENZYL ALCOHOL :UG/L :10 U :10 WSO8 1.2-DICHLOROBENZENE :UG/L :10 U :10	U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U	10 U 10 U 10 U 10 U 10 U	
WSO6 1.4-DICHLOROBENZENE UG/L 10 U 10 WSO7 BENZYL ALCOHOL UG/L 10 U 10 WSO8 1.2-DICHLOROBENZENE UG/L 10 U 10	U 10 U U 10 U U 10 U U 10 U U 10 U U 10 U	10 U 10 U 10 U	
WSO7 BENZYL ALCOHOL	U 10 U U 10 U U 10 U U 10 U	10 U 10 U	
WSOB 1.2-DICHLOROBENZENE UG/L 10 U 10	U 10 U U 10 U U 10 U	10 U	
	U 10 U	10 U	:
WSO9 2-METHYLPHENOL (O-CRESOL) :UG/L:10 U:10	U :10 U	-:	
	:	10 U .	
WS10 B1S(2-CHLOROISOPROPYL) ETHER :UG/L:10 U:10	U :10 U	-,	·
WS11 4-METHYLPHENOL (P-CRESOL) :UG/L:10 U:10		10 U	:
WS12 N-NITROSO-DIPROPYLAMINE : UG/L: 10 U:10	บ 10 บ	10 0	
WS13 HEXACHLOROETHANE :UG/L:10 U:10	U 10 U	10 U	
WS14 NITROBENZENE :UG/L:10 U:10	บ 10 บ	10 U	
WS15 ISOPHORONE : UG/L : 10 U : 10	U 10 U	10 U	
WS16 2-NITROPHENOL	U 10 U	10 U	
WS17 2 4-DIMETHYLPHENOL UG/L :10 U :10 U	U 10 U	10	
WS18 BENZOIC ACID :UG/L :50 U :50 L	บ :50 บ	:50 U :	
WS19 BIS(2-CHLOROETHYOXY) METHANE UG/L :10 U :10 U	U 10 U	10 U	
WS20 2.4-DICHLOROPHENOL UG/L :10 U :10 U	U 10 U	10 U	
WS21 1.2.4-TRICHLOROBENZENE UG/L :10 U :10 L	υ :10 υ	10 0	
WIZZ NAPHTHALENE UGZL 10 U 10 L	ນ :10 ປ	10 0	
WSZ3 4°CHLOROÁNH INI. UG/L :10 U :10 U	บ 10 เ	10 0	
WS24 HEXACHLOROBUTADICIJE UG/L : 10 U : 10 U	U 10 U	10 U	
WS25 4-CHLORO-3-METHYLPHENOL UG/L 10 U :10 U	J 10 U	10 0	
WS26 2-METHYLNAPHTHALENE UG/L 10 U :10 U	J 10 U	:10 U	
WS27 HEXACHLOROCYCLOPENTADIENE UG/L : 10 U : 10 U	J 10 U	10 U	
WS28 2.4.6-TRICHLOROPHENOL UG/L 10 U 10 U	J 10 U	10 U	

COMPOUND	UNITS	20	06		206D	4	207F	20	8F	209F
WS29 2.4.5-TRICHLOROPHENUL	: UG/L	: 25	U	: 25	U	: 25	U	25	υ :	
WS30 2-CHLORONAPHTHALENE	UG/L	10	Ü	10	U	10	U	10	U	
WS31 2-NITROANILINE (ORTHO NITROANILINE)	UG/L	25	U	:25	U	25	U	25	U	
WS32 DIMETHYLPHTHALATE	UG/L	10	U	10	U	10	U	10	U :	
WS33 ACENAPHTHYLENE	UG/L	10	U	:10	U	:10	U	: 10	U	
WS34 3-NITROANILINE	UG/L	25	U	:25	U	25	U	25	υ :	
WS35 ACENAPHTHENE	UG/L	10	U	:10	U	10	U	10	U	
WS36 2,4-DINITROPHENOL	UG/L	25	U	: 25	U	25	U	25	U :	
WS37 4-NITROPHENOL	UG/L	25	U	25	U	25	U	: 25	U	
W536 DIBENZOFURAN	UG/L	10	U	10	<u>-</u>	10	U	10	U	
WS39 2 4-DINITROTOLHENE	UG/L	10	U	10	บ	10	v	10	U	
WS40 2.6-DINITROTOLUENE	UG/L	10	U	10	U	10	U	10	U	
WS41 DIETHYLPHTHALATE	UG/L	10	Ū	10	Ü	10	U	10	U	
WS42 4-CHLOROPHENYL PHENYL ETHER	UG/L	10	U	10	U	10	U	10	U	
WS43 FLUORENE	UG/L	10	U	10	U	10	U	10	U	
WS44 4-NITROANILINE	UG/L	25	U	25	U	25	U	25	U :	
WS45 4,6-DINITRO-2-METHYLPHENOL	UG/L	25	U	25	U	25	บ	25	U	•
WS46 N NITROSODIPHENYLAMINE	UG/L :	10	U	10	U	10	U	10	U :	
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L	10	U	10	U	10	U	10	u :	
wide Hisachi okoßin, Lol	UG/L	10	U	10	()	10	11	10	U .	
wishs removem orderens	UG/L	2 5	Ü	25	U	25	U	25	U	
WOSO PHI NANTHRENE	:06/1	10	······································	10	U	10	U	10		
W351 -MINICACENE	UG/L	10	Ū	10	U	10	IJ ;	10	U	
WSS2 DI N BUTYL PHTHALAIL	UG/L :	10	Ū :	10	U :	10	U :	10		
WSSA FLUORANTHENE	UG/L	10	Ü	10	U	10	U	10	U :	:
WS54 PYRENE	UG/L	10	υ:	10	U	10	U	10	U	· · · · · · · · · · · · · · · · · · ·
	::-		:		:		:		:	:

	COMPOUND	UNIT	S	206	206	D	207F		208	F		209F
WS5!	5 BUTYL BENZYL PHTHALATE	UG/L	10	U	:10	 U	: 10	u	: 10	 U	-: :	
WS56	3.3'-DICHLOROBENZIDINE	UG/L	10	U	10	U	10	U	10	U	-:	
WS5	BENZO(A)ANTHRACENE	UG/L	: 10	U	10	U	:10	U	10	U	-:	
WS58	BIS(2-ETHYLHEXYL)PHTHALATE	UG/L	10	U	10	U	10	U	10	U	-:	
WS59	CHRYSENE	UG/L	10	U	10	U	10	U	10	U	- :	
WS60	DI-N-OCTYL PHTHALATE	UG/L	:10	V	10	Ü	10	U	10	U	- :	
WS61	BENZO(B)FLUORANTHENE	UG/L	10	U	10	U	10	U	10	Ū	- :	
W562	BENZO(K)FLUORANTHENE	UG/L	10	U	10	U	10	U	10	U	:	
W563	BENZO(A)PYRENE	UG/L	10	U	10	U	10	U	10	U		
W564	INDENO(1,2,3-(D)PYKENE	UG/L	10	 U	10	U	10	IJ	10	IJ		
w565	DIBENZO(A,H)ANTHRACENE	UG/L	10	U	: 10	U	10	U	10	U		:
W566	BENZO(G.H,I)PERYLENE	UG/L	10	U	10	U	10	U	10	U		
WS67	CARBAZOLE	UG/L	10	U	10	U	10	U	10	U		
WT09	CYANIDE, TOTAL	MG/L	0.017	7 U	0.010	U	0.010	U	0.010	U		
WV03	CHL OROME THANE	UG/L	10	U	10	U	10	U	10	U	10	y :
WV04	BROMOMETHANE	UG/L	10	U	10	U	10	U	10	U	10	U :
WV05	VINYL CHLORIDE	UG/L	10	V	10	U	10	U	10	u u	10	ป
wv06	CHI OROE THANE	UG/L	10	U	10	U	10	U	10	U	10	U
WV07	METHYLENE CHLORIDE	UG/L	: 10	U	10	U	10	U	10	υ	10	U
WV()8	1.1-DICHEOROFTHEME	:06/1	10	U	10	; U	10	11	10	11	10	II .
MA03	1, 1-DICHLOROETHANE	UG/L	10	U	10	U	10	Ü	10	Ü	10	U
W710	1.2-DICHLOROETHENE TOTAL	UG/I	10	U :	10	U	10	IJ.	10	IJ	10	U
WVII	CHILOROFORM	UG/L	: 10	Ü	10	U :	10	U :	10	Ü	10	U
WV12	1.2-DICHLOROETHANE	UG/L	10	U	10	U .	10	U	10	U	. 10	U .
WV13	1.1.1-TRICHLORGETHANE	UG/L	10	U	10	U	10	U	10	Į)	. 10	U .
WV14	CARBON TETRACHLORIDE	UG/L	10	U	10	U	10	U	10	U	10	0
				:		:		:				

COMPOUND	UNITS	206		206D		2076	:	208F		209	Э F
BROMOD I CHLOROME THANE	: UG/L	: 10	U	10	U	: 10	U	10	U	:10	U
1.2-DICHLOROPROPANE	UG/L	10	U	:10	U	10	U	10	ับ	10	U
BENZENE	UG/L	10	U	: 10	U	10	U	10	U	: 10	U
TRICHLOROETHENE	UG/L	10	. U	: 10	U	10	U	10	U	10	U
CIS-1,3-DICHLOROPROPENE	UG/L	10	U	:10	U	:10	U	10	u	10	U
DIBROMOCHLOROMETHANE	UG/L	10	U	: 10	U	10	U	10	U	10	Ü
1,1,2-TRICHLOROETHANE	UG/L	: 10	ีบ	:10	U	10	U	10	บ	10	U
BROMOFORM	UG/L	10	U	10	U	: 10	U	10	U	10	Ū
TETRACHLOROETHENE	UG/L	10	U	10	Ū	10	U	10	υ	10	υ
TOLUENE	UG/L	: 10	v	: 10	Ū	10	U	10	t	10	U
1,1,2,2-TETRACHLOROETHANE	:UG/L	: 10	U	: 10	U	10	IJ	10	U	10	U
CHLOROBENZENE	UG/L	10	U	: 10	Ū	10	U	10	U	: 10	U
ETHYL BENZENE	UG/L	10	U	: 10	U	10	U	10	บ	: 10	U
ACETONE	UG/L	10	U	: 10	U	10	U	10	U	: 10	U
CARBON DISULFIDE	UG/L	10	U	10	U	10	U	10	U	10	U
2-BUTANONE	:UG/L	10	U	10	U	10	U	10	U	10	υ :
VINYL ACETATE	UG/L	10	U	10	U	10	U	10	U	10	΄υ :
2-HEXANONE	UG/L	10	U .	10	บ	10	U	: 10	U	10	U .
4-METHYL-2-PENTANONE	UG/L	10	U	10	U	10	U	10	11	10	U
STYRENE	UG/L	10	u	10	U :	10	u	10	11	10	U
XYLENES TOTAL	:06/L	10	U	10	U	10	U	10	H	10	U
TRANS-1.3-DICHLOROPROPENE	UG/L	10	บ	10	บ :	10	(1	10	0	10	0
HYDROCARBONS, TOTAL PETROLEUM	MG/L		:		: :	0.5	IJ				• •
SAMPLE NUMBER	. NA	206	:	206	: :	207		203		209	
ACTIVITY CODE	NA	DSX44	: :	DSX44	: :	DSX44		DSX-I-I		DSX44	:
	COMPOUND BROMODICHLOROMETHANE 1.2-DICHLOROPROPANE BENZENE TRICHLOROETHENE CIS-1.3-DICHLOROPROPENE DIBROMOCHLOROMETHANE 1.1.2-TRICHLOROETHANE BROMOFORM TETRACHLOROETHENE 1.1.2.2-TETRACHLOROETHANE CHLOROBENZENE ETHYL BENZENE ACETONE CARBON DISULFIDE 2-BUTANONE VINYL ACETATE 2-HEXANONE 4-METHYL-2-PENTANONE STYRENE XYLENES FOLAL TRANS-1.3-DICHLOROPROPENE HYDROCARBONS, TOTAL PETROLEUM SAMPLE NUMBER ACCITVITY CODE	BROMODICHLOROMETHANE 1.2-DICHLOROPROPANE UG/L BENZENE UG/L TRICHLOROETHENE UG/L CIS-1.3-DICHLOROPROPENE UG/L DIBROMOCHLOROMETHANE UG/L 1.1.2-TRICHLOROETHANE UG/L BROMOFORM UG/L TETRACHLOROETHENE UG/L 1.1.2.2-TETRACHLOROETHANE UG/L CHLOROBENZENE UG/L ETHYL BENZENE UG/L ACETONE CARBON DISULFIDE UG/L VINYL ACETATE UG/L	BROMODICHLOROMETHANE	BROMODICHLOROMETHANE UG/L 10 U 1.2-DICHLOROPROPANE UG/L 10 U BENZENE UG/L 10 U TRICHLOROETHENE UG/L 10 U TRICHLOROPROPENE UG/L 10 U DIBROMOCHLOROMETHANE UG/L 10 U 1.1.2-TRICHLOROETHANE UG/L 10 U \$\$ROMOFORM UG/L 10 U TETRACHLOROETHENE UG/L 10 U TOLUENE UG/L 10 U 1.1.2.2-TETRACHLOROETHANE UG/L 10 U CHLOROBENZENE UG/L 10 U ACETONE UG/L 10 U ACETONE UG/L 10 U CARBON DISULFIDE UG/L 10 U 2-BUTANONE UG/L 10 U VINYL ACETATE UG/L 10 U VINYL ACETATE UG/L 10 U 4-METHYL-2-PENTANONE UG/L 10 U XYLENES 101AL UG/L 10 U XYLENES 101AL PETROLEUM MG/L SAMPIE NUMBER NA 206	BROMODICHLOROMETHANE	BROMODICHLOROMETHANE	BROMODICHLOROMETHANE	BROMODICHLOROMETHANE	BROMODICHLOROMETHANE UG/L 10 U 10 U	BROMODICHLOROMETHANE UG/L 10 U 10 U 10 U 10 U 10 U 10 U 10 U 10	BROMODICHLOROMETHANE

ANALYSIS REQUEST DETAIL REPORT ACTIVITY: 1-DSX44

		c	OMPOUND	UNIT	s 30	01	301	D	302		303		304	
WFO1	WATER TEMP	· •		· c	: 7. O		.7.0		6.0		:6.0		. 3 0	
WFOS	PH. FIELD			SU	8.15		8.15		7.87		7 . 27		: 7.05	, -
WF 10	CONDUCTIVITY (FIE	LD)	: UMHO	5:440		440		450		360		350	
WMO 1	SILVER	BY	ICAP	UG/L	10	U	10	U	10	Ū	10	U	: 10	U
WMO2	ALUMINUM	BY	ICAP	UG/L	5100	J	4800	J	5000	J	28000		10000	J
WMO3	ARSENIC	ВУ	ICAP	UG/L	10	U	10	U	10	U	10	U	:10	U
WMO4	BARIUM	BY	ICAP	UG/L	200	U	200	U	200	U	280		200	U
WMO5	BERYLLIUM	BY	ICAP	UG/L	5.0	U	5.0	U	5.0	U	5.0	U	5 0	U
WM06	CADMIUM	ВУ	ICAP	UG/L	5.0	U	5.0	U	5.0	U	5.0	U	5 0	U .
WMO7	UBALT	Вγ	ICAP	UG/L	50	U	50	U	50	U	50	U	50	ii ii
WM08	CHROMIUM	BY	ICAP	UG/L	10	U	10	U	10	U	31	IJ	: 14	
WMO9	COPPER	BY	ICAP	UG/L	25	U	25	U	25	U	25	U	25	U :
WM10	IRON	BY	ICAP	UG/L	5500		5300		5500		27000		11000	
WM 1 1	MANGANESE	BY	ICAP :	UG/L	250		250		250		600		450	:
WM12	MOLYBDENUM	BY	ICAP	UG/L	NA	0	NA	0	NA	0	NA	0	MA	0
WM13	NICKEL	BY	ICAP	UG/L	40	U	40	U	40	Ü	40	U	40	Ü:
WM14	LEAD	BY	ICAP	:UG/L	7 0		7.2		9.7		24	U	18	•
WM15	ANT IMONY	BY	ICAP	:UG/L	60	U	60	U	60	U	60	U	:60	U
WM16	SELENIUM	BY	ICAP	:UG/L	5.0	U	5.0	U	5.0	U	NA	I	5 0	U
wai 7	T COMPOS	Б <i>і</i>	10 AP	UG/L	: NA	0	NA NA	0	NA	()	NA	ű.	· MA	O.
WMTo	FHGI L LUM	BY	ICaP	UG/L	10	Ü	10	U	10	1)	MA	t	10	Ü
WM19	MAHADTHM	BY	ICAP	UG/L	50	U	50	U :	50	11	62		r.0	II
WiA2O	21M	Ü'n	ICAP	UG/L	28	U	26	U	33	. U :	89	U	e' t	
WH21	CALCIUM, TOTAL	BY	ICAP	.MG/L	53		54		54	:	51	11	. 58	
WI122	MAGNESTUM, TOTA	L B	Y ICAP	:MG/L	18		18	:	18	·	19	Ū	23	
WM23	SODIUM, TOTAL B	ΥÏ	CAP	MG/L	36	:	36	:	36		30	U:	19	:

	COMI	POUN	D	UNITS	•	301	;	301D	302		30	3		304
WM24	POTASSIUM, TOTAL BY	ICA	P	MG/L	6 2		6.1		6.2		:8.8	U	:5.2	·
WM35	SILVER.DISSOLVED	BY	ICAP	UG/L	10	U	10	U	10	U	10	บ	10	U
WM36	ALUMINUM.DISSOLVED	ВУ	ICAP	UG/L	200	U	200	U	200	U	200	U	200	V
WM37	ARSENIC, DISSOLVED	BY	ICAP	UG/L	10	U	10	U	10	· U	10	U	10	U :
WM38	BARIUM.DISSOLVED	BY	ICAP	UG/L	200	Ü	200	· U	200	U	200	U	200	U
WM39	BERYLLIUM, DISSOLVED	ВУ	ICAP	UG/L	5 0	U	5.0	U	5.0	U	5.0	U	5.0	Ū :
WM40	CADMIUM, DISSOLVED	BY	ICAP	UG/L	5.0	U	5.0	U	5.0	U	5.0	U	5.0	U
WM41	COBALT.DISSOLVED	BY	ICAP	UG/L	50	, υ	50	· U	50	U	50	U	50	U :
WM42	CHROMIUM, DISSOLVED	BY	ICAP	UG/L	10	U	10	U	10	U	10	U	10	U :
WM43	COPPER DISSOLVED	ВУ	ICAP	UG/L	25	U	25	U	25	U	25	U	25	U
WM14	IRON, DISSOLVED	BY	ICAP	UG/L	100	U	100	U	100	U	100	U	100	U :
WM45 I	MANGANESE, DISSOLVED	BY	ICAP	UG/L	17		15	U	15	U	15	U	15	U
WM46 I	MOLYBDENUM, DISSOLVED	BY	ICAP	UG/L	NA	0	NA	0	NA	0	NA	0	NA .	0
WM47 N	NICKEL . DISSOLVED	BY	ICAP,	UG/L	40	V	40	U	40	U	40	U	40	U
WM48 I	FAD DISSOLVED	BY	ICAP	UG/L	3.0	U	3.3	J	3.0	U	3 0	U	3 0	ıi
WM49 A	ANTIMONY, DISSOLVED	BY	ICAP	UG/L	60	U	60	U :	60	U	60	U	60	V :
WM50 S	SELENIUM.DISSOLVED	BY	ICAP	UG/L	5.0	U	5.0	U	5.0	U	11	J	5 0	41
WM51 T	ITANIUM.DISSOLVED	BY	ICAP	UG/L	NA	0	NA	0	NA	0	NA	0	: NA	0
WM52 T	HALL TUM. DISSOLVED	BY	ICAP	UG/L	10	U	10	U	10	U	10	U	10	U
WIA W	Mian (M. 04550) vrn	ftY	ICAP	.0671	50	()	50	U :	50	U	50	li	40	1)
WMF-4	THE DESCRIPTION	bit	ICAP.	0071	20	U	20	U	20	U	20	U	20	U
WMS5 C	ALCIUM.DISSOLPED	EV	ECVB	MG/L:	52		52	:	54		46 .		53	
WIASO M	AGRESTOM, DISSOLVED	ВУ	1CAP	MG/L	17		17	· · · · · · · · · · · · · · · · · · ·	18		15		20	
WAS7 S	ODTUM, DISSOLVED	D/	ICAP	MG/L	36		37		38		30		. 19	
WM to 1	OTA551UM.DESOLVED	87	ICAP	MG/L:	5.3		5.2	: (5.4		5.2		5.0	U .
W501_P	HENOI			UG/L	10	U	10	U :	10	U	10	U	10	U

COMPOUND	UNITS	301		3	301D	302	2	30	3	3	304
WSO3 BIS(2-CHLOROETHYL) ETHER	UG/L	: 10	U	10	U	10	U	10	u	10	U
WSO4 2-CHLOROPHENOL	UG/L	10	U	10	U	10	U	10	U	10	U
WSO5 1,3-DICHLOROBENZENE	UG/L	10	U	10	U	10	U	10	 U	10	U
WSO6 1.4-DICHLOROBENZENE	UG/L	10	U	10	U	10	U	:10	U	10	Ü
WSO7 BENZYL ALCOHOL	UG/L	10	U	10	U	10	U	10	U	10	U
WSO8 1,2-DICHLOROBENZENE	UG/L	10	U	10	Ü	10	U	10	U	10	U
WSO9 2-METHYLPHENOL (O-CRESOL)	UG/L	10	U	10	U	10	U	10	U	10	U
WS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/L	10	U	10	U	10	U	10	U	: 10	U :
WS11 4-METHYLPHENOL (P-CRESOL)	UG/L	10	U	10	U	10	U	10	U	10	U
WS12 N-NITROSO-DIPRUPYLAMINE	UG/L	10	U	10	U	10	U	10	U	10	U
WS13 HEXACHLOROETHANE	UG/L	10	U	10	U	10	U	10	U	10	U ·
WS14 NITROBENZENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS15 ISOPHORONE	UG/L	10	U	10	v	10	U	10	U	10	U
WS16 2-NITROPHENOL	UG/L	10	U	10	U	10	U	10	V	10	U
WS17 2.4-DIMETHYLPHENOL	UG/L	10	U	10	U	10	U	10	U	10	.11
WS18 BENZOIC ACID	UG/L	50	U	50	U	50	v	50	U	50	Ü :
WS19 BIS(2-CHLOROETHYOXY) METHANE	UG/L	10	U	10	U	10	U	10	U	10	· U
WS20 2.4-DICHLOROPHENOL	UG/L	10	U	10	U	10	U	10	U	10	U :
WS21 1.2.4-TRICHLOROBENZENE	UG/L	10	U	10	U	10	U	10	U	10	U
WWW NAPHIHALENE	UG/L :	10	U	10	U	10	11	10	U	10	U
Wind distill UROANTI TNE	UG/L	10	U	10	U	10	U	10	II.	10	Ü
W321 HEXACHI OROBUTADTEHI	UG/I	10	U	10	Ü :	10	U	10	IJ	10	U
W525 4- CHLORO - 3-METHYLPHENOL	UG/L :	10	U	10	Ü:	10	U :	10	U .	10	U
WS2G 2 METHYLNAPHTHALENE	:UG/L :1	10	U	10	U	10	U	10	U :	10	U ,
WS27 HEXACHLOROCYCLOPENTADIENE	UG/L 1	10	U	10	Ü	10	U	10	Ü	10	ii
WS28 2.4.6-TRICHLOROPHENOL	UG/L 1	10	U :	10	Ū	10	Ü	10	U :	10	U :
	::-			:	:		:				:

COMPOUND	UNIT	S	301		301 D		302		303		304
WS29 2,4,5-TRICHLOROPHENOL	: UG/L	25	U	: 25	U	25	U	25	u	25	U
WS30 2-CHLORONAPHTHALENE	UG/L	10	U	10	U	10	U	10	U	10	V
WS31 2-NITROANILINE (ORTHO NITROANILINE)	UG/L	25	U	25	U	25	U	25	U	25	Ü
WS32 DIMETHYLPHTHALATE	UG/L	10	U	:10	บ	10	U	10	U	10	U
WS33 ACENAPHTHYLENE	UG/L	10	U	:10	Ū	10	U	10	U	10	U
WS34 3-NITROANILINE	UG/L	25	U	25	บ	25	U	25	U	25	U
WS35 ACENAPHTHENE	UG/L	10	U	:10	U	10	U	10	U	10	U
WS36 2.4-DINITROPHENOL	UG/L	25	U	25	U	25	U	25	U	25	U :
WS37 4-NITROPHENOL	UG/L	25	U	25	U	: 25	U	25	U	25	U
WS38 DIBENZOFURAN	UG/L	10	U	10	U	10	U	10	U	10	U ;
WS39 2,4-DINITROTOLUENE	UG/L	10	U	: 10	Ü	10	U	10	บ	10	U
WS40 2.6-DINITROTOLUENE	:UG/L	10	V	10	U	10	U	10	U	10	U
WS41 DIETHYLPHTHALATE	UG/L	10	U	10	U	10	U	10	U	10	Ü
WS42 4-CHLOROPHENYL PHENYL ETHER	UG/L	10	Ū	10	U	10	U	10	U	10	Ü
WS43 FLIIORENE	:UG/L	10	บ	10	Ü	10	U	10	U	10	U
WS44 4-NITROANILINE	UG/L	25	U	25	U	25	U	25	U	25	Ü
WS45 4.6-DINITRO-2-METHYLPHENOL	UG/L	25	U	25	U	25	U	25	U	25	,U
WS46 N-NITROSODIPHENYLAMINE	UG/L	10	บ	10	U	10	U	10	U	10	U
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L	10	U	10	Ü	10	U	10	U .	10	U
WS48 HEXACHLOROBEN /FILL	067L	10	ļ)	10	U	10	U	10	11	10	U :
WS49 PENTACHLOROPHENOL	UG/L	25	U	25	U	25	U	25	U ·	25	u i
WSSO PHENANTHRENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS51 ANTHRACENE	UG/L	10	U	10	U	10	U	10	U :	10	U
WS52 DI-H-BUTYL PHTHALATE	UG/L	10	U	10	V	10	U	10	U .	10	U
WS53 FLUORANTHENE	:UG/L	10	U	10	U	10	U .	10	U :	10	U .
WS54 PYRENE	UG/L	10	IJ	10	U	10	U :	10	U	10	U
					;		-				•

	•										
COMPOUND	UN1T:	S 3	101	3010		302		303		304	1
WS55 BUTYL BENZYL PHTHALATE	UG/L	10	U	10	U	10	 U	10	U	10	U .
WS56 3.3'-DICHLOROBENZIDINE	UG/L	10	U	10	U	10	U	10	U	10	U
WS57_BENZO(A)ANTHRACENE	: UG/L	10	U	10	υ	10	U	10	U	10	U
WS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/L	10	U	:10	U	10	U	10	U	10	U
WS59 CHRYSENE	UG/L	10	U	10	U	10	U	10	U	10	U :
WS60 DI-N-OCTYL PHTHALATE	UG/L	10	U	10	U	10	U	10	U	10	U
WS61 BENZO(B)FLUORANTHENE	UG/L	10	U	10	U	10	U	10	U	10	U
WS62 BENZO(K)FLUORANTHENE	UG/L	10	U	10	U	10 -	U	10	U	10	U .
WS63 BENZO(A)PYRENE	UG/L	:10	U	10	U	: 10	U	10	U	10	U .
WS64 INDENO(1.2.3-CD)PYRENE	UG/L	10	U	10	U	10	U	10	U	10	U :
WS65 DIBENZO(A.H)ANTHRACENE	UG/L	10	U	10	U	10	U	10		10	
WS66 BENZO(G.H.I)PERYLENE	:UG/L	10	U	10	U	10	U	10	U	10	U :
WS67 CARBAZOLE	:UG/L	10	U	10	U	10	U	10	U	10	U :
WTO9 CYANIDE, TOTAL	MG/L	0.010	U	0.010	U	0.010	U	0.010	U	0 010	U :
WVO3 CHLOROMETHANE	UG/L	10	U	10	U	10	H	10	ti.	10	H .
WVO4 BROMOMETHANE	:UG/L	10	U	10	U	10	U	10	U	10	U
WVO5 VINYL CHLORIDE	UG/L	10	Ü	10	U	10	U	10	U	10	Ú
WVO6 CHLOROETHANE	UG/L	10	U	10	U	10	U	10	U	10	U :
WVO7 METHYLENE CHLORIDE	UG/L	10	U	10	U	10	U	10	U	10	U
WVO8 1 1: DICHLOROFTHENE	UG/1	10	U	10	U	10	11	10	II.	10	U
WVO9 1.1 DECHLOROE HIADE	: UG/L	10	U	10	U	10	u :	10	0	10	
WY10 1 2 DICHLOROETHENE. TOTAL	: UG/1.	10	U	10	Ü	10	U.	10	U	10	U
WV11 CHLOROFORM	UG/L	10	U	10	U	10	u .	10	u	10	ti
WV12 1.2 DICHLOROETHANE	UG/L	10	U	10	Ü	10	U.	10	U	10	Ü.
WV13 1.1.1 TRICHLORGETHANE	UG/L	10	U	10	Ü .	10	U :	10	U.	10	Ü
WV14 CARBON TETRACHLORIDE	UG/L	10	Ū	10	Ü	10	U	10	U	10	U .

	COMPOUND	UNITS	5	301		301D	3	102	30	3	304	4
WV15	BROMODICHLOROMETHANE	: UG/L	:10	U	10		: 10	 U	10	U	. 10	 i
WV16	1.2-DICHLOROPROPANE	UG/L	10	U	10	U	10	U	10	U	10	U
WV17	BENZENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV19	TRICHLOROETHENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV20	CIS-1,3-DICHLOROPROPENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV21	DI BROMOCHLOROME THANE	UG/L	10	V	10	U	10	U	10	U	10	U
WV22	1.1.2-TRICHLOROETHANE	UG/L	10	U	10	V	10	U	:10	U	10	U
WV24	BROMOFORM	UG/L	10	U	10	U	10	U	10	U	10	U
WV25	TETRACHLOROETHENE	UG/L	10	ຸ ບ	10	บ	10	บ	10	υ	10	บ
WV26	TOLUENE	: UG/L	10	U	10	U	10	U	10	 U	10	U U
WV27	1.1.2.2-TETRACHLOROETHANE	UG/L	10	U	10	U	10	U	10	U	10	11
WV28	CHLOROBENZENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV29	ETHYL BENZENE	UG/L	10	U	10	U	10	U	10	U	10	U
WV30	ACETONE	UG/L	10	V	11	U	10	Ü	10	U	10	U
WV3,1	CARBON DISULFIDE	UG/L	10	U	10	U	10	U	10	11	10	U
WV32	2-BUTANONE	UG/L	10	U	10	U	10	V	10	U	10	U
MA33	VINYL ACETATE	UG/L	10	U	10	U	10	U	10	U	10	บ
WV34	2-HEXANONE	UG/L	10	U	10	U	10	V	10	U	10	U
WV35	4-METHYL-2-PENTANONE	UG/L	10	U	10	U	10	U	10	บ	10.	U
MA III	STYRENE	UG/L	10	()	10		10	U	10	U	10	11
wy.i/	AYLENES TOTAL	U6/L	10	U	10	U	10	II	10	Ü	10	10
W.,40	TRANG-1.3-DICHLOROPROPENE	UG/L	10	U	10	U	10	IJ	10	Ü	10	0.1
WV36	HYDROCARBONS, TOTAL PETROLEUM	MG/L	NA	0	0.5	Ū	0.5	U	0.5	U	5 5	U :
ZZ01	SAMPLE NUMBER	NA	301		301		302		303	·	304	
2702	ACTIVITY CODE	NA .	DSX44		DSX44		D5X44		DSX44	 	D5X44	

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-DSX44

	COMPOUND	UNITS	304D	401	4011	D	402		40	3
SMO1 SILVER	BY ICAP	MG/KG:	: 2 . 6	·	J:2.6	ี	:2.6	U	2.6	U
SMO2 ALUMINUM	BY ICAP	MG/KG:	550)	6800		12000		10000	
SMO3 ARSENIC	BY ICAP	MG/KG:	3.7		4.0		8.4		7.1	
SMO4 BARIUM	BY ICAP	MG/KG:	140		140		160		160	
SMO5 BERYLLIUM	BY ICAP	MG/KG:	1.3	(1.3	U	1.3	U	1.3	li .
SMO6 CADMIUM	BY ICAP	MG/KG:	1.3	(1.3	U	1.3	U	1.3	ບ
SMO7 COBALT	BY ICAP	MG/KG	13	(13	U	13	U	13	U
SMO8 CHROMIUM	BY ICAP	MG/KG:	8.7		9.0		16		:12	
SMO9 COPPER	BY ICAP	MG/KG	9.1		8.2		26		23	
SM10 IRÓN	BY ICAP	MG/KG:	9800		10000		19000		14000	
SM11 MANGANESE	BY ICAP	MG/KG:	250		300		320		250	J
SM12 MOLYBDENUM	BY ICAP	MG/KG	NA NA	0	NA	0	NA	0	NA	0
SM13 NICKEL	BY ICAP	MG/KG	10		11		18		: 16	
SM14 LEAD	BY ICAP	MG/KG:	30		13	J	36	J	31	J
SM15 ANTIMONY	BY ICAP	MG/KG	15	U	15	U	16	U	16	'n
SM16 SELENIUM	BY ICAP	MG/KG:	:1.3	บ	1.3	บ	2.0	J	1.3	U
SM17 TITANIUM	BY ICAP	MG/KĠ	NA NA	0	NA	0	NA	0	NA	-0
SM18 THALLIUM	BY ICAP	MG/KG	:2.6	U	2.6	U	2.6	Ü	2.6	U
SM19 VANADIUM	BY TCAP	MG/KG	: 15		17		27		25	
SM2O 2116	Br (CBP	MG/KG:	: 35	J	36	J	77	J	64	
SMST CALCIUM	By DAP	MG/KG:	6400		6300		6500	• • • •	5/00	
SM22 MACHESTUM	BY TOAP	MG/KG:	2700		2900	 :	3000		3600	
SN23 SODIUM	Br TCAP	MG/KG:	: 1300	U	1300	U :	1300	U	1300	U
SM24 POTASSIUM	BY ICAP	:MG/KG	1300	U	1300	U	1700		1600	
SSOT PHENOL		UG/KG:	420	U	410	U .	820	U	1800	U .
5502 CARBAZOLE		UG/KG:	420	U	410	U	820	ย	1800	Ū

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COMPOUND	UNITS	304D	401		40	1D	402	:	40	3
SSO3 BIS(2-CHLOROETHYL) ETHER	UG/KG:		420	U	410	-	820		1800	U
SSQ4 2-CHLOROPHENOL	UG/KG:		420	U	410	U	820	U	1800	U
SSOS 1,3-DICHLOROBENZENE	UG/KG:		420	U	410	U	820	 U	1800	 U
SSOG 1,4-DICHLOROBENZENE	UG/KG		420	U	410	U	820	U	1800	U
SSO7 BENZYL ALCOHOL	UG/KG:		NA .	• 0	NA .	0	NA	0	: NA	0
SSOB 1.2-DICHLOROBENZENE	UG/KG:		420	U	410	U	820	U	1800	U
SSO9 2-METHYLPHENOL (O-CRESOL)	UG/KG:		420	U	410	U	820	U	1800	U
SS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/KG:		420	U	410	U	820	U	1800	U
SS11 4-METHYLPHENOL (P-CRESOL)	UG/KG:		420	U	410	U	820	U	1800	ť
SS12 N-NITROSO-DIPROPYLAMINE	UG/KG:		420	U	410	U	820	l!	1800	ti
SS13 HEXACHLOROETHANE	UG/KG:		420	U	410	U	820	U	1800	U
SS14 NITROBENZENE	UG/KG:		420	U	410	U	820	IJ	1800	U
SS15 ISOPHORONE	UG/KG:		420	U	410	U	820	U	1800	U
SS16 2-NITROPHENOL	UG/KG		420	U	410	U	820	Ü	1800	U
SS17 2.4-DIMETHYLPHENOL	UG/KG:		420	U	410	บ	820	U	1800	U
SS18 BENZOIC ACID	UG/KG:		: NA	0	NA .	0	NA	O	NA	0
SS19 BIS(2-CHLOROETHYOXY) METHANE	UG/KG		420	U	410	U	820	U	1800	าเ
SS20 2.4-DICHLOROPHENOL	UG/KG:		420	U	410	U	820	U	1800	U
SS21 1.2.4-TRICHLOROBENZENE	UG/KG:		420	U	410	U	820	U	1800	U
SS22 NAPHTHALENE	UGZKG:		420	U	410	U	820	U	1800	U
5523 4-CHLOROANTETNE	:UG/KG:		420	U	410	u :	820	U	1800	ii .
SS24 HEXACHI OROBUTADIENE	UG/KG		: 120	IJ	: 410	u :	820 .	и.	1200	U i
5525 4 CHLORO-3-METHYLPHEHOL	UG/KG:		420	U .	410	U .	820	U.	1800	U .
SS26 2 METHYLNAPHTHALENE	.UG/KG:		420	บ	410	U	820	ü,	1800	u .
SS27 HEXACHLOROGYCLOPENTADIENE	UG/KG:		420	U	410	U	820	Ü	1800	Ü
SS28 2.4.6-TRICHLOROPHENOL	UG/KG		420	U	410	U	820	<u> </u>	1800	<u></u>

ANALYSIS REQUEST DETAIL REPORT

ACTIVITY: 1-DSX44

COMPOUND	UNITS	304D	401		401D	402		403	
SS29 2.4.5-TRICHLOROPHENOL	:UG/KG:		: 1000	U	: 990	υ : 2000	 U	4300	U
SS30 2-CHLORONAPHTHALENE	UG/KG:		420	U	410	820	U	1800	U
SS31 2-NITROANILINE	UG/KG:		1000	Ü	990	2000	U	4300	U
SS32 DIMETHYLPHTHALATE	UG/KG:		420	U	410	820	U	1800	Ü
SS33 ACENAPHTHYLENE	UG/KG:		420	U	410	820	U	1800	U
SS34 3-NITROANILINE	UG/KG		1000	U	990	2000	U	4300	Ü
SS35 ACENAPHTHENE	UG/KG:		420	U	410 L	820	U	1800	U
SS36 2.4-DINITROPHENOL	UG/KG:		1000	U	990 U	2000	U	4300	U
SS37 4-NITROPHENOL	UG/KG:		1000	U	. 990 U	2000	U	4300	U
SS38 DIBENZOFURAN	UG/KG:		420	U	410 U	820	H	: 1800	Ü
SS39 2,4-DINITROTOLUENE	UG/KG:		420	U	410 U	820	U	1800	11 :
SS40 2.6-DINITROTOLUENE	:UG/KG:		420	U	410 U	820	ี ป	1800	U
SS41 DIETHYLPHTHALATE	UG/KG:		420	U	410 U	820	U	1800	U
SS42 4-CHLOROPHENYL PHENYL ETHER	UG/KG		420	U	410 U	820	U	1800	t:
SS43 FLUORENE	UG/KG:		420	u	410 U	820	IJ	1800	Ų.
SS44 4-NITROANILINE	UG/KG		1000	U	990 U	2000	υ	4300	U
SS45 4,6-DINITRO-2-METHYLPHENOL	UG/KG		1000	U	990 U	2000	U	4300	-11
SS46 N-NITROSODIPHENYLAMINE	UG/KG		420	U	410 U	:820	U .	1800	U
SS47 4-BROMOPHENYL PHENYL ETHER	UG/KG:		420	U	410 U	820	U	1800	IJ
SSUK HEXACHLOROBENZENE	UG/KG		420	u	410 0	820	H	1800	11
SSAPE PENEACHLORUPHENOL	UG/NG:		1000	U	990 U	2000	11	4300	ıi
SS50 PHENANTHRENE	UG/KG:		420	U :	410 "	\$2900		4400	•
SSST ANTHRACENE	UG/KG:		420	U	410 U	820	IJ.	1300	IJ
SS52 DI-H-BUTYL PHTHALATE	UG/KG:		420	U :	410 U	. 620	U	1300	Ü
SS53 FLUGRANTHENE	UG/KG		420	u 🥞	750	5000		5100	فند
SS54 PYRENE	UG/KG		420	U	60	6400	J	9000	

COMPOUND	UNITS	304D	401	40	1D	402		403	3
SSSS BUTYL BENZYL PHTHALATE	UG/KG	420	U	410	บ	:820	U	1800	U
SS56 3.3'-DICHLOROBENZIDINE	UG/KG	420	U	410	U	820	U	1800	U
SS57 BENZO(A)ANTHRACENE	UG/KG:	420	U	60	و المالية المالية المالية المالية المالية المالية المالية المالية المالية المالية المالية المالية المالية الم	3500	ala:	4200	
SS58 BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG:	420	Ū	410	U	820	U	: 1800	U
SS59 CHRYSENE	:UG/KG:	420	U	70		3900		4900	
SS60 DI-N-OCTYL PHTHALATE	: UG/KG:	: 420	U	410	U	820	U	1800	U STORES
SS61 BENZO(B)FLUORANTHENE	UG/KG:	: 420	U	:410	U	3600		3600	
SS62 BENZO(K)FLUORANTHENE	UG/KG:	420	U	410	U	\$900	, ,	3100	
SS63 BENZO(A)PYRENE	: UG/KG:	: 420	U	4 30 .		2600		5600	
SS64 INDENÚCL 2 3-CDIPYRENE	UG/KG:	420	U	:410	U	: 3000		2900	
SS65 DIBENZO(A, H)ANTHRACENE	UG/KG:	420	U	410	U	820	IJ	1800	U :
SS66 BENZO(G.H.I)PERYLENE	UG/KG:	420	U	410	U	\$500		4100	
STO9 CYANIDE	MG/KG	6.4	V	6.4	บ	6	ू पूर	6.5	ນີ້
SVO3 CHLOROMETHANE	UG/KG	13	U	13	U	13	U	14	U
SVO4 BROMOMETHANE	UG/KG:	:13	U	:13	U	13	()	14	1)
SVOS VINYL CHLORIDE	:UG/KG:	:13	U	: 13	U	13	U	14	Ū:
SVO6 CHLOROETHANE	UG/KG:	:13	U	13	U	13	U	14	ข
SVO7 METHYLENE CHLORIDE	UG/KG:	13	U	: 15	U	13	U	14	U
SV08 1.1-DICHLOROETHYLENE	UG/KG:	13	U	13	U	13	U	14	U :
170 C. L. J. (GCHLOROF HAME	UG/KG:	13	U	:13	l)	13	u	1.1	11
Notice Thousand 2-DICHLOROF HIVLENE	UG/KG:	: 13	U	13	11	13	H	14	Ü
1 11 CH GROFORM	UG/KG:	:13	U	: 13	U .	13	 U	14	Ü
1942 1 2 DICHLOROL HARIE	:UG/KG:	:13	U	13	U :	13	U	14	U
19713 T. L. F. TRICHLOROL THANK	UG/FG.	13	U	13	U	13	U.	14	U .
SOLIA CARBON LETRACHLORIDE	UG/KG:	13	U	13	U	13	Ü	14	Ü
SV15 BROMODICHLOROMETHANE	UG/KG:	:13	U	13	U	13	Ü	14	U :
					:	+	·· :		:

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	COMPOUND	UNITS	304D	401	401D	402	403
WMO2 ALUMINUM	BY ICAP	:UG/L :970)O J	÷:	-:	·:	- <u>,</u>
WMO3 ARSENIC	BY ICAP	UG/L 10	U	:	-:	: :	:
WMO4 BARIUM	BY ICAP	UG/L 200) U	:	-:	: :	:
WMO5 BERYLLIUM	BY ICAP	UG/L 5.0) U	:	:	:	:
WMO6 CADMIUM	BY ICAP	UG/L 5.0	U	:	-:	:	:
WMO7 COBALT	BY ICAP	UG/L 50	Ü				
WMO8 CHROMIUM	BY ICAP	UG/L 12					
WMO9 COPPER	BY ICAP	UG/L 25	U				
WM10 IRON	BY ICAP	UG/L 110	00				
WM11 MANGANESE	BY ICAP	UG/L 430					·
WM12 MOLYBDENUM	BY ICAP	UG/L NA	0	:	:		:
WM13 NICKEL	BY ICAP	UG/L :40	U	:		:	
WM14 LEAD	BY ICAP	UG/L 15				:	:
WM15 ANTIMONY	BY ICAP	UG/L 60	Ū	:		:	
WM16 SELENIUM	BY ICAP	UG/L 5 0	U	:		:	
WM17 TITANIUM	BY ICAP	UG/L NA	0			:	:
WM18 THALLIUM	BY ICAP	UG/L 10	U				
WM19 VANADIUM	BY ICAP	UG/L 50	U				
WM20 ZINC	BY ICAP	UG/L 46					
WM21 CALCIUM TOTA	AL BY ICAF	MG/L 57					
WM22 MAGNESTUM. To	MAL BY LLAP	MG/L 22					•
WM23 SODIUM, TOTAL	BY TEAP	MG/I. 19					·
WM24 POTASSIUM, TO	OTAL BY ICAP	MG/L :5 1					•
WM35 SILVER, DISSOL	VED BY ICAP	UG/L 10	U				
WM36 ALUMINUM.DISS	OLVED BY ICAP	:UG/L :200	U				
WM37 ARSENIC.DISSO	DLVED BY ICAP	UG/L 10	U				
····		, :+	:		:		

ANALYSIS REQUEST DETAIL REPORT ACTIVITY: 1-DSX44

COMP	POUND	UNITS	304D		401	401D	402	403
WM38 BARIUM, DISSOLVED	BY ICAP	UG/L	200	U	:	-	-;	
WM39 BERYLLIUM, DISSOLVED	BY ICAP	UG/L	5.0	Ū	:	: :	:	
WM40 CADMIUM.DISSOLVED	BY ICAP	UG/L	5.0	U	:	:	:	
WM41 COBALT, DISSOLVED	BY ICAP	: UG/L	:50	U	:	:	· · · · · · · · · · · · · · · · · · ·	
WM42 CHROMIUM.DISSOLVED	BY ICAP	UG/L	10	U	:	:	: :	
WM43 COPPER, DISSOLVED	BY ICAP	UG/L	25	U	:	:	:	:
WM44 IRON.DISSOLVED	BY ICAP	UG/L	100	U		:	:	:
WM45 MANGANESE, DISSOLVED	BY ICAP	UG/L	15	U	:	:	:	
WM46 MOLYBDENUM.DISSOLVED	BY ICAP	UG/L	NA	0	:	:		
WM47 NICKEL DISSOLVED	BY ICAP	UG/L	40	U		:		
WM48 LEAD.DISSOLVED	BY ICAP	UG/L	3.0	U		:		· · · · · · · · · · · · · · · · · · ·
WM49 ANTIMONY, DISSOLVED	BY ICAP	UG/L	60	U		:	:	:
WM50 SELENIUM, DISSOLVED	BY ICAP	UG/L	5.0	Ū		:	:	:
WM51 TITANIUM.DISSOLVED	BY ICAP:	UG/L	NA	0		:	:	:
WM52 THALLIUM.DISSOLVED	BY ICAP	UG/L	10	U :		:	:	
WM53 VANADIUM.DISSOLVED	BY ICAP	UG/L	50	υ:		:	:	
WM54 ZINC.DISSOLVED	BY ICAP	UG/L	20	U		:		
WM55 CALCIUM.DISSOLVED	BY ICAP	MG/L	55	:	*****	:	:	· · · · · · · · · · · · · · · · · · ·
WM56 MAGNESTUM. DISSOLVED	BY ICAP	MG/L	21	:		:	:	·
WMSZ SODTOM DESSOLVED	BY TOAP	MG/L	20	:				
Willias FOTASSTOM DISSOLVED	BY ICAP	MG/L	5.0	U :			: • • • • • • • • • • • • • • • • • • •	
W1G1 PHICHOL		.UG/L	10	U :				
w 03 Bluc2 CHEOROLTHYL) ET	HER	UG/L	10	U :	:	:	;	
WTO4 2 CHLORGPHENOL		UG/L	10	U		·	:	
W505 1.3 DICHLOROBENZEHE		UG/L	10	: U :	:		· · · · · · · · · · · · · · · · · · ·	
WHO6 1.4 DICHLOROBENZEHE		UG/L	10	:		· · · · · · · · · · · · · · · · · · ·		
		-::		:	:			,

•	ANALYSIS REQUE	ST DETAIL REPORT	TAIL REPORT ACTIVITY: 1-DSX44			
COMPOUND	UNITS	304D	401	401D	402	403
WSO7 BENZYL ALCOHOL	UG/L 10	v :			-;	
WSOB 1.2-DICHLOROBENZENE	UG/L 10	U			:	:
WSO9 2-METHYLPHENOL (O-CRESOL)	UG/L 10	U			: :	:
WS10 BIS(2-CHLOROISOPROPYL) ETHER	UG/L 10	U			:	
WS11 4-METHYLPHENOL (P-CRESOL)	UG/L 10	U			:	;
WS12 N-NITROSO-DIPROPYLAMINE	UG/L 10	Ū			:	:
WS13 HEXACHLOROETHANE	UG/L 10	U			:	-:
WS14 NITROBENZENE	UG/L 10	U	:		:	
WS15 ISOPHORONE	UG/L 10	U			:	:
W516 2-NTTROPHENOL	UG/L 10	U	:		:	
WS17 2.4-DIMETHYLPHENOL	UG/L 10	U			•	
W518 BENZOIC ACID	UG/L 50	U			:	:
WS19 BIS(2-CHLOROETHYOXY) METHANE	UG/L :10	Ü				:
WS20 2.4-DICHLOROPHENOL ;	UG/L 10	U				
WS21 1 2 4-TRICHEOROBENZENE	UG/L 10	Ü			:	
WS22 NAPHTHALENE	UG/L 10	v				
WS23 4~CHLOROANILINE	UG/L 10	Ū	:			
WS24 HEXACHLOROBUTADIENE	UG/L :10	U				:
WS25_4-CHLORO-3-METHYLPHENOL	UG/L 10	U			:	
Wile of State of the Herbert Miles	U6/L 10	U :	:			
Willer HE COUNTY ON THE PROPERTY OF THE PARTY OF THE	UG/L 10	U :	:		:	
wi.25 2 1.6- IRTCHLOROLULION	UG/L 10	U	:		•	
W. 20-2, 4 % TRICHEOROPHENOL	UG/L .25	U			, -	
WESO 2 OH OKOHAPITHALLING	UG/L 10	U	:		:	••
wish 2 MITROAHIITHE CORTHO NITROANILINE)	UG/L 25	U			The second secon	
WS32 DIMETHYLPHTHALATE	UG/L 10	Ü	:			

COMPOUND	UNITS	304D	401	401D	402	403
WS33 ACENAPHTHYLENE .	UG/L :10	U :		-; 	:	·
WS34 3-NITROANILINE	UG/L 25	Ü		:	:	:
WS35 ACENAPHTHENE	UG/L 10	U				
WS36 2.4-DINITROPHENOL	UG/L 25	U		-:	:	:
WS37 4-NITROPHENOL	UG/L 25	U		-:		
WS38 DIBENZOFURAN	UG/L 10	Ü				
WS39 2.4-DINITROTOLUENE	UG/L 10	U		-:		
WS40 2,6-DINITROTOLUENE	UG/L 10	U		-:		:
WS41 DIETHYLPHTHALATE	UG/L 10	U		:	:	:
WS42 4-CHI OROPHENYL PHENYL ETHER	UG/L 10	Ü		:	:	
WS43 FLUORENE	UG/L 10	υ		:	:	
WS44 4-NITROANILINE	UG/L 25	U				:
WS45 4.6-DINITRO-2-METHYLPHENOL	UG/L 25	U			:	
WS46 N-NITROSODIPHENYLAMINE	UG/L 10	U				
WS47 4-BROMOPHENYL PHENYL ETHER	UG/L 10	U :		:		
WS48 HEXACHLOROBENZENE	UG/L :10	U :		:		•
WS49 PENTACHLOROPHENOL	UG/L 25	U		:		
WS50 PHENANTHRENE	UG/L :10	U :		:		:
WS51 ANTHRACENE	UG/L :10	U :		:		:
WS2 PINGUTYL PHIHALAIF	UG/L 10	Ü		:		·
W-53 FI UORANTHENE	UG/L 10	U		, 	· · · · · · · · · · · · · · · · · · ·	
WS54 PYRENE	:UG/L :10	U		:	·	
WSSS BUTYL BENZYL PHTHALATE	. UG/L : 10	V :		:		
WOWG 3.31-DICHLOROBENZIDINE	UG/L :10	U :				
W 657 BEN (O (A) ANTHRACENE	UG/L :10	U :		:		
WSSR BIS(2-ETHYLHEXYL)PHIHALATE	UG/L 10	י ט :		:	:	

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VALIDATED DATA

ANALYSIS REQUEST DETAIL REPORT

:DSX44

2202 ACTIVITY CODE

:DSx44

VALIDATED DATA ACTIVITY: 1-DSX44 401 401D 402 403

D5444

D5X44

COMPOUND UNITS 304D :UG/L :10 U: WV20 CIS-1.3-DICHLOROPROPENE U: WV21 DIBROMOCHLOROMETHANE :UG/L :10 WV22 1,1,2~TRICHLOROETHANE :UG/L :10 WV24 BROMOFORM :UG/L :10 WV25 TETRACHLOROETHENE :UG/L :10 U : WV26 TOLUENE :UG/L :10 WV27 1,1,2,2-TETRACHLOROETHANE :UG/L :10 U: WV28 CHLOROBENZENE :UG/L :10 **u** : WV29 ETHYL BENZENE : UG/L : 10 WV30 ACETONE :UG/L :10 WV31 CARBON DISULFIDE : UG/L : 10 U : : UG/L : 10 U : WV32 2-BUTANONE WV33 VINYL ACETATE :UG/L :10 U: WV34 2-HEXANONE :UG/L :10 U WV35 4-METHYL-2-PENTANONE :UG/L :10 U **WV36 STYRENE** :UG/L :10 U: WV37 XYLENES, TOTAL :UG/L :10 U WV40 TRANS-1,3-DICHLOROPROPENE U: :UG/L :10 WV86 HYDROCARBONS. TOTAL PETROLEUM MG/L :0 5 U : 2201 SAMPLE NUMBER 402 : NA : 304 : 401

: DSX44

COMPOUND UNITS 501 SSO1 PHENOL : UG/KG: 450000 SSO3 BIS(2-CHLOROETHYL) ETHER : UG/KG: 450000 U : SSO4 2-CHLOROPHENOL U : : UG/KG: 450000 SSO5 1.3-DICHLOROBENZENE : UG/KG: 450000 U: SSO6 1.4-DICHLOROBENZENE : UG/KG: 450000 U : SSO7 BENZYL ALCOHOL : UG/KG: 450000 U: SSOB 1,2-DICHLOROBENZENE : UG/KG: 450000 U: SSO9 2-METHYLPHENOL (O-CRESOL) : UG/KG : 450000 U SS10 BIS(2-CHLOROISOPROPYL) ETHER UG/KG: 450000 U U SS11 4-METHYLPHENOL (P-CRESOL) : UG/KG: 450000 SS12 N-NITROSO-DIPROPYLAMINE : UG/KG: 450000 U: SS13 HEXACHLOROETHANE : UG/KG: 450000 SS14 NITROBENZENE : UG/KG: 450000 U: SS15 ISOPHORONE : UG/KG: 450000 U: SS16 2-NTTROPHENOL UG/KG: 450000 U SS17 2.4-DIMETHYLPHENOL U : UG/KG: 450000 SS18 BENZOIC ACID UG/KG: 2300000 SS19 BIS(2-CHLOROETHYOXY) METHANE UG/KG: 450000 SS20 2.4-DICHLOROPHENOL U : UG/KG: 450000 \$521 1.2.4-TRICHLOROBEN/ENF UG/KG:450000 U SS22 NAPHTHALENE : UG/KG **# 00000** 5323 4 CHLOROANTLINE : UG/KG: 450000 U : 5524 HEXACHLOROBUTADIENE : UG/KG: 450000 SS25 4-CHLORO-3-METHYLPHENOL : UG/KG: 450000 : UG/KG**T100000Q** SS26 2-METHYLNAPHTHALENE SS27 HEXACHLOROCYCLOPENTADIENE : UG/KG: 450000

ACTIVITY: 1-DSX44

ANALYSIS REQUEST DETAIL REPORT

COMPOUND	UNITS 501				
SS28 2.4.6-TRICHLOROPHENOL	: UG/KG: 450000	U		:	· ; •
SS29 2.4.5-TRICHLOROPHENOL	UG/KG: 2300000		:	:	
SS30 2-CHLORONAPHTHALENE	UG/KG: 450000	U	:	:	
SS31 2-NITROANILINE	UG/KG: 2300000	U	:		
SS32 DIMETHYLPHTHALATE	:UG/KG:450000	U			
SS33 ACENAPHTHYLENE	:UG/KG: 450000	U			
SS34 3-NITROANILINE	UG/KG: 2300000	U		:	
SS35 ACENAPHTHENE	UG/KG: 620000			:	
SS36 2.4-DINITROPHENOL	UG/KG: 2300000	Ū		:	
SS37 4-NITROPHENOL	UG/KG: 2300000	U	:		
SS38 DIBENZOFURAN	UG/KG: 450000	V	:	·	
SS39 2.4-DINITROTOLUENE	UG/KG: 450000	V		· · · · · · · · · · · · · · · · · · ·	
SS40 2.6-DINITROTOLUENE	UG/KG: 450000	U		: :	:
SS41 DIETHYLPHTHALATE	UG/KG: 450000	U	:		:
SS42 4-CHLOROPHENYL PHENYL ETHER	UG/KG: 450000	U	:	· · ·	:
SS43 FLUORENE	:UG/KG: 820000 (表表表。	i get			
SS44 4-NITROANILINE	: UG/KG: 2300000	Ū			: :
SS45 4.6-DINITRO-2-METHYLPHENOL	:UG/KG: 2300000	U			:
SS46 N-NITROSODIPHENYLAMINE	:UG/KG: 450000	U			· ;
SS47 4-BROMOPHENYL PHENYL ETHER	: UG/NG: 450000	U :			
5548 HEXACHLOROBENZENE	: UG/KG: 450000	U :			:
SS49 PENTACHLOROPHENOL	UG/KG: 2300000	U		:	
SS50 PHEMANTHRENE	UG/KG: 4300000				:
SSS1 ANTHRACENE	UG/KG: \$10000	· () : :			
SS52 DI-N-BUTYL PHTHALATE	UG/KG: 450000	U			
SS53 FILUORANTHENE	:UG/KG:				

ACTIVITY: 1-DSX44

COMPOUND	UNITS	501				
SS54 PYRENE	UG/KG	300000	:	:		::
SSSS BUTYL BENZYL PHTHALATE	UG/KG: 45	50000 U	:	:		
SS56 3.3'-DICHLOROBENZIDINE	UG/KG:90	00000 U		:		
SS57 BENZO(A)ANTHRACENE	UG/KG 🛕6	00000			:	
SS58 BIS(2-ETHYLHEXYL)PHTHALATE	:UG/KG:45				:	:
SS59 CHRYSENE	UG/KG: 27	00000			:	
SS60 DI-N-OCTYL PHTHALATE	UG/KG:45	0000 U	:	:	:	:
SS61 BENZO(B)FLUORANTHENE	UG/KG: 45	0000 U	:	· :	· :	
SS62 BENZO(K)FLUORANTHENE	: UG/KG: 45	0000 U	:	:	:	
SS63 BENZO(A)PYRENE	UG/KG: 27	ي نيد کړي 0000	:	:	· : ·	
SS64 INDENO(1,2,3-CD)PYRENE	: UG/KG: 45	0000 U		:		
SS65 DIBENZO(A,H)ANTHRACENE	UG/KG: 45	0000 U	· :	· ·		:
SS66 BENZO(G,H,I)PERYLENE	UG/KG: 450	0000 U	· :	· · ·		· · · · · · · · · · · · · · · · · · ·
SVO3 CHLOROMETHANE	UG/KG:500	00 U		· · ·	·	:
SVO4 BROMOMETHANE	UG/KG: 500	00 U	· ·		· ·	
SV05 VINYL CHLORIDE	UG/KG:500	00 U	, ` , , , , , , , , , , , , , , , , , ,			
SVO6 CHLOROETHANE	UG/KG:500	00 U				•
SVO7 METHYLENE CHLORIDE	:UG/KG:500	00 U				
SVO8 1.1-DICHLOROETHYLENE	UG/KG: 250	00 U				
SVO9 1,1-DTCHLORGETHANE	:UG/KG: 250)O U				
SVIO TRANS-1.2-DICHLOROETHYLENE	: UG/KG: 250	00 U				
SV11 CHLOROFORM	UG/KG: 250	00 U				
SV12 1.2-DICHLOROETHANE	UG/KG: 250)O U				
SV13 1.1.1-TRICHLOROETHANE	UG/KG: 250	00 U			1	
SV14 CARBON TETRACHLORIDE	UG/KG: 250	00 U				
SV15 BROMODICHLOROMETHANE	UG/KG: 250	ο υ				:

COMPOUND	UNITS	501				ı	
SV16 1.2-DICHLOROPROPANE	UG/KG: 2500) U	:	:	:	· , 	٠.
SV17 BENZENE	UG/KG: 2100		!	:	:	:	:
SV18 TRANS-1.3-DICHLOROPROPENE	UG/KG 2500			:	:	:	:
SV19 TRICHLOROETHYLENE	UG/KG: 2500) U	:	:		:	:
SV20 CIS-1.3-DICHLOROPROPENE	UG/KG: 2500) U		:			:
SV21 DIBROMOCHLOROMETHANE	UG/KG: 2500	Ü					
SV22 1.1.2-TRICHLOROETHANE	UG/KG: 2500	U			:	:	:
SV24 BROMOFORM	UG/KG: 2500	U	:	:	:	:	:
SV25 1.1.2.2-TETRACHLOROETHENE	UG/KG: 2500	U	:	:	: :	· :	:
SV26 FOLUENE	UG/KG:8400	0	· ·		: :	· : :	:
SV27 1.1.2.2-TETRACHLOROETHANE	UG/KG 2500	U	· : ·	:	· ! '	· · ·	:
SV28 CHLOROBENZENE	UG/KG: 2500	U	: : : !	:	· · ·	· :	:
SV29 ETHYL BENZENE	UG/KG: 3800	0	· : ·	:	, , , ,	· · ·	:
SV30 ACETONE	UG/KG: 5000	U		· :		· · ·	:
SV31 CARBON DISULFIDE	UG/KG: 2500	U		:	, , , ,	, ; ; ;	
SV32 2-BUTANONE	UG/KG: 5000	U	· · ·	•		· · ·	:
SV33 VINYL ACETATE	UG/KG: 5000	U	· ·	•			:
SV34 2-HEXANONE	UG/KG:5000	U	· ·	•		· · ·	
SV35 4-METHYL-2-PENTANONE	UG/KG:5000	U		· · ·		· · ·	
SMB6 STYRENE	UG/KG: 2500	U :		: :			
SWW XALLMES TOTAL	UG/KG: 19600	00					
SV54 HYDROCARBONS, TOTAL PETROLEUM	MG/KG: ATTAC	HMENT					
2201 SAMPLE NUMBER	NA 501						
ZZO2 ACTIVITY CODE	NA DSX44	!		· ·			

ACTIVITY DSX44 LACLEDE COAL GAS

THE PROJECT LEADER SHOULD CIRCLE ONE - STORET, AIRS, OR ARCHIVE.

CIRCLE ONE:

STORET

AIRS

ARCHIVE

DATA APPROVED BY LABO FOR TRANSMISSION TO PROJECT LEADER ON 06/12/91 09:01:15 BY

FOR ACTIVITY: DSX44

SPFD

06/12/91 16:34:32

* FINAL REPORT

FY: 91 ACTIVITY: DSX44 DESCRIPTION: LACLEDE COAL GAS LOCATION: MISSOURI

STATUS: ACTIVE TYPE: SAMPLING - CONTRACT LAB ANALYSIS PROJECT: A34

LABO DUE DATE IS 5/14/91. REPORT DUE DATE IS 5/28/91.

INSPECTION DATE: 3/14/91 ALL SAMPLES RECEIVED DATE: 03/15/91

ALL DAIA APPROVED BY LABO DATE: 06/12/91 FINAL REPORT TRANSMITTED DATE: 06/21/91

EXPECTED LABO TURNAROUND TIME IS 60 DAYS EXPECTED REPORT TURNAROUND TIME IS 75 DAYS

ACTUAL LABO TURNAROUND TIME IS 89 DAYS ACTUAL REPORT TURNAROUND TIME IS 99 DAYS

SAMP.	QCC	M	DESCRIPTION	SAMPLE STATUS		CITY	STATE	AIRS/ STORET LOC NO	BEG. DATE	BEG. Time	END. DATE	END. Time
001	-	S	BACKGROUND SAMPLE-CSI-A	1		LOUIS	MISSOURI		03/07/91	15:36	/ /	
002		Š	BACKGROUND SAMPLE-CSI-B	i		Louis	MISSOURI		03/07/91	11:00	1 1	•
003		Š	BACKGROUND SAMPLE-CSI-C	i	4 ST.	LOUIS	MISSOURI		03/07/91	16:45	11	i
004		Š	BACKGROUND SAMPLE-CSI-D	i	4 ST.	LOUIS	MISSOURI		03/07/91	16:55	11	:
005		S	BACKGROUND SAMPLE-CSI-E	1	4 ŠŤ.	LOUIS	MISSOURI		03/07/91	17:15	7 7	:
006		S	BORING LOCAL. #6-B-06-CS2	1	4 ST.	LOUIS	MISSOURI		03/08/91	09:20	11	
007		S	BORING LOCAL #6-R-06-CS2	1	4 ST.	LOUIS	MISSOURI		03/08/91	09:35	11	:
800		S	R-06-CS3 SAMPLE LOCATION	1	4 ST.	LOUIS	MISSOURI		03/08/91	11:07	1 1	:
009		S	B-14-CS4 SAMPLE LOCATION	1	4 ST.	LOUIS	MISSOURI		03/08/91	13:50	11	:
010		W	B14-CS4 SAMPLE LOCATION	1	4 ST.	LOUIS	MISSOURI		03/08/91	14:20	/ /	:
011		5	B-07-CS6 SAMPLE LOCATION	1	4 ST.	LOUIS	MISSOURI		03/08/91	17:35	/ /	:
012			R13-USZ SAMPLE LOCATION	1	4 ST.	roniz	MISSOURI		03/09/91	09.24	/ /	:
013		•	B13 (S/ SAMPLE LOCATION	1	4 SŢ.	roniz	MISSOURI		03/09/91	09 45	/ /	
joj		•	BORING LOCATION-01/MOUND STR POWER	1	3 51	LOUIS	MISSOURI		03/04/91	13 05	/ /	
102		ž.	BORING LOCATION 02	1	3 51.	LOUIS	MISSOURI		03/02/91	14 20	/ /	
103		ς.	SURFACE SOIL AT EAST CENTRALOF TANK	ļ	3 ST.	LOUIS	MISSOURI		03/04/91	15.30	/ /	
104			BORING LOCATION 05	1		LOUIS	MISSOURI		03/04/91	17.25	/ /	
105		?	BETWEEN TANKS 7-8/BORING 07	!	3 ST.	LOUIS	MISSOURI		03/05/91	13:50	/, /,	:
106		2	BORING LOCATION-15				MISSOURI		03/06/91	14:10	1. 1.	
107		ج	BORING LOCATION -16	!	3 ST.	roniz	MISSOURI		03/06/91	15: 15	/ /	
108		5	BORING LOCATION-17		3 ST.	roniz	MISSOURI		03/06/91	15:45	1, 1,	:
109 110		5	BORING LOCATION-18	!		roniz	MISSOURI		03/06/91	16:53	', ',	:
111		Š	R19-SAMPLE LOCATION	!		LOUIS	MISSOURI		03/14/91	07:45	', ',	:
112		5	R2O-SAMPLE LOCATION	!		LOUIS	MISSOURI		03/14/91	07:30	', ',	:
112		3	B21-SAMPLE LOCATION	1	3 ST.	LOUIS	MISSOURI		03/13/91	17:30	/ /	•

APPENDIX E

Field Sheets and Chain of Custody Records

RAFT		S. ENVI		TAL PR		ION /	AGENCY				_	4445	
y: 91 4													
STIVITY			-		P 2 0	JECT	NUM:	434	PT:	LAT:	SITU!	DE:	
AMPLE E DCATION ASE/BAT TORET/S	CES: _ N: ST. ICH/SMI SAROAD	LOUIS	i - A		LAS:		8 E G E N E	6: 03 0:	DATE 5/61/ //	91 <u>15</u>	:36	=ROM R EAST: NORTH: DOWN:	
ALYSIS CASS CASS CASS CASS CASS	Q	PRES ICE(NON(ICE(NON)	}		55 54 54 5109	VOL.	ATTLES ALSO (s.		COL	D VAI	P ilil	uti •
CTMENT	S: '	SEG BA	e CK		_		aFedy)	, L	'S #	· L			,

Policy considerated for

AMPLE COLLECTED BY : PECK/MCLALL

U. ENVIRONMEN	S. ENVIRONMENT TAL SERVICES D	AL PROTECTION IV. 25 FUNSTO	AGENCY, REG N RD. KANSAS	ION VII	115
: 71 ACTNO:	DSX44 SAMNG:	002 9CC: _ MED	IA: SOIL P	L: S P F D	
CATION:	LACLEDE COAL	MO PROJECT	NUM: A34 P		
MPLE DES: CATION: ST. SE/BATCH/SM DRET/SARGAD	LOUIS 0:/_	MO LAB:	DA BEG: 03/6 END:/_	11807 TIME FO H 191 /1:00 E	ROM REF PT AST: ORTH: OWN: 7-11/F+
ALYSIS REQU NTAINER ASS ASS ASS ASS	ORESERVATIV ICED NONE ICED	SS (SEM	ATILES ALS	BY COLD VAP	At-to
MMENTS:	JEE BO	2P	1 NS + SAI	PETY LOC	た本し

Muci concentration.

CAPI	•	CIECO PHEEL		•	
ENVIRG	U.S. ENVIRONMENTAL NMENTAL SERVICES DIS	_			KS 66115
': 91 AC	TNG: DSX44 SAMMG: 00)3 2CC: _ ME	DIA: SOIL	PL: S P	F D
	DES: LACLEDE COAL GA		T NUM: A3	REF LATI 4 PT: LONG	
ICATION:	S: CS -C ST. LOUIS H/SHO:/ :CN DADS	MO LAB:	AFG:	กรี/ดา/91 //	ME FROM REF PT : \(\sigma \) EAST: NORTH:
HALYSIS	REQUESTED:				
INTAINER	PRESERVATIVE	MGP NA	ME		
_ASS	ICED	55 (58	MIVOLATIL	ES je to toto	-
_ASS	NONE		LATILES		
-ASS	ICED_	-SM VOIE		_	1.1171
ASS-	NONE	- STC9 iC		-	
-ASS	ICED	SH34 ME		SY COLD	VAP /

MMERTS:

SEE BORING + SAF ETY LOG#2 Background

ficial. Concentation

AMPLE COLLECTED EY: PECK/mcCALL

ENVIRONM	U.S. ENVIRONMENTA MENTAL SERVICES DI					
': 91 ACTN	NG: DSX44 SAMNO: 0	04 QCC: _	MEDIA: SOI	L PL: S	PPD	*****
	S: LACLEDE CJAL G		JECT NUM: 4	REF LA'		
MPLE DES: CATION: S SE/BATCH/ ORET/SARG	ST. LOUIST /SMO:/_	MO LAB:	BEG:	*5DATE 63/61/91 / - /01/_	TIME FROM L:SEAST: NORTH	REF PT
ALYSIS RENTAINER ASS ASS ASS	PRESERVATIVE ICED NONE ICED	/55 5Y /5H /5T09	SEMIVOLATI VOLATILES	- extinctor	 •	
MMENTS:	SEE BOR BAC	ING & S	SAFETY OUND	L06#	2	

Mild Christophytion

MPLE COLLECTED 34 : PELK / Malal

	.S. ENVIRONMENTAL NTAL SERVICES DIV				66115
: 01 ACTNO	: 35X44 SAMNC: 90	15 4CC: _ MED	IA: SOIL	PL: S P F D	
CATION: 5	; LACLEDE COAL G	40 PROJECT	NUM: 434	REF LATITUD PT: LONGITU	
DRET/SARDA	(5)-E LOUIS MC: // D NO:	MO LAB:	BEG: 0	DATE TIME 3/81/91 /Z:/S 107/:_	EAST:
ALYSIS REG NTAINER NSS ASS ASS ASS	PRESERVATIVE ICED NONE ICED NONE ICED NONE ICED	'SS SEM SV VOL MET STO9CYA	HIVOLATILE ATILES ALS_P	Sy COLD VA	- Atti
MMENTS:	SEE Boring BALKGRE	_			
		- V	rued	conent	whi

MPLE COLLECTED 34 : PECK/MacALL

		FIELD SHEET TAL PROTECTION AGENCY, REGION VII DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115	
91 ACTNO:	DSX44 SAMNO:	006 RCC: _ MEDIA: SOIL PL: S P F D	
	LAGLEDE COAL	GAS REF LATITUDE: MO PROJECT NUM: A34 PT: LONGITUDE:	
APLE DES:	C:	MO BEG: 03/07/91 07: 20 EAST: LAB: END: // NORTH: DOWN: 3-5/0	<u>e</u>
ALYSIS REQU	ESTED: PRESERVATI	VE MGP NAME /	
155	ICED	SS SEMIVOLATILES Wellie	
155	NONE	SV VOLATILES (SM METALS: (STOY CYANIDE)	
122	ICED	SH HETALS.	
128	ICFD	SM34 MERCURY 84-COLD VAP 8	
		N.O. I	
		no Hg:	
1MENTS:			
	Comment of the second	Location 76 continuous.	
	Liung		
	U.,		
	sample		
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	Su l	onny 9 refer log. #2	
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		. 110	
		hud concentration	
		MUST CONTRACT	
		1 Color Color Lichton	

APLE COLLECTED BY : POCK () Nich and

	. ENVIRONMENTAL AL SERVICES DIV				56115
: 91 ACTNO:	DSX44 SAMNG: 00	7 QCC: _ MEDI	A: SOIL	PL: S P F D	
	LACLEDE COAL GAS		NUM: A34	REF LATITUDI PT: LONGITU	
MPLE DES: _/ CATION: ST. SE/BATCH/SMC DRET/SAROAD	LOUIS 1: NO:	MO LAB:	8EG: 03	DATE TIME \$104/91 14:15 104/ _:	FROM REF PT EAST: NORTH: DOWN: P-13-ft
ALYSIS REQUE	STED: PRESERVATIVE	MGP NAM	E	r /	
	ICED	SEM SEM		S. C. C. C. C. C.	-
ASS	NONE	SV VOL	ATTLES	del	
ASS ASS			ALS;		_
AS-S			CURY	BY COLO VA	FAO Hage
MMENTS:		•			
	Bound !	oration	#6 c	ontinor-	⇒
	sample?	チン		4)
	Lu l	our &	sofeti	log #0	<
		./1-			
		Mil	V. Co	M.	

	.S. ENVIRONMENTAL NTAL SERVICES DIV	PROTECT				
: 91 ACTNO	: DSX44 SAMNC: 90	8 OCC: _	MEDIA: SOIL	. PL: S P	F D	
TIVITY DES	: LACLEDE COAL GA		JECT NUM: A3	REF LATI		
MPLE DES: CCATION: ST SE/BATCH/S DRET/SARDA	<u>19-06- (53</u> SMO:/_/ NO NO:/	MO _ LAB:	BEG:	03/91/91 //	ME FROM R (:27 EAST: : NORTH: DOWN:	2-12 fz
IALYSIS REG	NESTED:					
INTAINER		MGP	NAME			
.ASS	ICED	5 S	SEMIVOLATIL	-ES		
.ASS	NONE	SV	VOLATILES			
.ASS	ICED	SM				
.ASS	NONE	ST09	CYANIDE			_
.ASS	ICED	SH34	MERCURY	BY COL	LYAPAUL	eti.

IMMENTS:

lee bourg i rafety log #2

V/D med-conc

AMPLE COLLECTED BY : FRCT 1216 COLL

U.S. ENVIRONMENTAL PROTECTION AGENCY/ REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 : P1 ACTNO: DSX44 SAMNO: 009 QCC: _ MEDIA: SOIL PL: S P F D TIVITY DES: LACLEDE COAL GAS REF LATITUDE: CATION: ______MO PROJECT NUM: A34 PT: LONGITUDE: MPLE SES: 3.14 - 654 お CATE TIME FROM BEG: 03/94791 (公:52) EAST: FROM REF PT MO CATION: ST. LOUIS SE/BATCH/SMO: ____/_ LAB: ___ END: __/20/___:_ NORTH: DRET/SARDAD NO: ALYSIS REQUESTED: MGP NTAINER PRESERVATIVE NAME ASS ICED SS SEMIVOLATILES S¥ ASS NONE VOLATILES ASS ICED SM METALS STO9 CYANIDE ASS NONE ASS ICED MMENTS: Lee boung & safety logbook #2 V/13 mid. Conc.

FIELD SHEET

AFT

ENVIRONMEN'	S. ENVIRONMENTAL Tal services div.	. 25 FUNSTO	AGENCY, REG In RD. Kansas	CITY, KS á	6115
r: 91 ACTNG:	DSX44 SAMNG: 019	O QCC: _ MED	IA: SOIL P	L: S P F D	
TIVITY DES:	LACLEDE COAL GAS	S MO PROJECT	R Num: 434	REF LATITUDE PT: LONGITUS	E:
AMPLE DES: _ DCATION: ST. ASE/EATCH/SM TORET/SAROAD	パゲ・ 〜 〜 〜 〜 〜 LOUIS O:/_/	CM LAB:	デオ DA BEG: 03/4 END:/	ATE TIME 97/91 / <u>9</u> :30	FROM REF PT EAST: NORTH: DOWN: 22-74-ft
NALYSIS PEGU ONTAINER LASS LASS LASS LASS LASS	ESTED: PRESERVATIVE	NGP NAM S SEP SV VOL	ME MIVOLATILES LATILES TALS ANTOF		· delate
OMMENTS:	Su bou oil 4	water	fety de	vegbook um es	#2 neintration
			i Darn 1- Hun	•	JATR

	S. ENVIRONMENTAL Tal services div.				
: 91 ACTNB:	DSX44 SAMNG:)11	CC: _ MEDIA:	SOIL PL: S F	P F D	***
CATTON:	LACLEDE COAL GAS	MO SPOLECT MIN	REF LAT : A34 PT: LOT	ICTTUDE.	
MPLE DES: CATION: ST. SE/24TCH/SM ORET/S4ROAD	Λ-07- & Δ΄ LOUIS O:/_ NO:	MO B	GDATE 1 EG: C3/91/91/ NO: _/_/_	TIME =ROM RE 7:15 EAST: NORTH: COWN:	F PT
ALYSIS REQU	ESTED: PRESERVATIVE ICED NONE	MGP NAME SS SEMIVOL SV VOLATIL SM METALS	ATILES Es	o var del	·
MMENTS:	See bour	y Safe	ty logio.	x ≠3	
		V/B	med, Con	ı.	

FIELD SHEET

MPLE COLLECTED BY : 1978 / CE-AREA

AFT U. ENVIRONMEN	S. ENVIRONMENTAL	FIELD SHEET PROTECTION AGENCY, REGION VII 25 FUNSTON RD. KANSAS CITY, KS 66115	
7: 91 ACTNO:	DSX44 SAMNO: 01	2 QCC: _ MEDIA: SOIL PL: S P F D	
CATION:		MO PROJECT NUM: A34 PT: LONGITUDE:	
AMPLE DES: DCATION: ST. ASE/BATCH/SM CORET/SAROAD	K/3- ∠57 LOUIS 0:/ No:	### ADATE TIME FROM R MO BEG: 03/01/91 09:2/ EAST: LAB: = NO: 10/1 = NORTH: DOWN:	EF PT
22A. 22A.	ESTED: PRESERVATIVE ICED NONE ICED NONE ICED NONE ICED	MGP NAME SS SEMIVOLATILES SV VOLATILES SM METALS STD9 CYANIDE SM34 MERCURY BY COLD VAP	Lest:
CTMENTS:	See bor	ing + safety logbook #	5
		V/B med Cone	

FIELD SHEET RAFT U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 1: 31 ACTNO: DSX44 SAMNO: 013 QCC: _ MEDIA: SDIL PL: S P F D AMPLE DES: SITE NO A DATE TIME FROM REF PT BEG: 03/01/91 09:45 EAST: 4SE/BATCH/SHO: ____/_/__ LAB: __ END: _______ NORTH: TORET/SARDAD NO: ____ IALYSIS REQUESTED: INTAINER PRESERVATIVE MGF NAME ICED SS .ASS SEMIVOLATILES _ASS SV VOLATILES _ASS ICED SM METALS -ASS NONE 90T2 CYANIDE BY COLD VAP () (Little ICED .ASS SM34 MERCURY IMMENTS: Au boing & sufety logbook #3

V/B Med. Cone

AMPLE COLLECTED BY : FROM Stropes

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS. 66115
: 91 ACTNO: 05X44 SAMNO: 101 OCC: _ MEDIA: SOIL PL: S P F 0
TIVITY DES: LACLEDE COAL GAS REF LATITUDE: CATION: _B \(\text{O} \)
MPLE DES: OPECATION: BD MO PROJECT NUM: A34 PT: LONGITUDE: OPECATION: ST. LOUIS MO BEG: 03/94/91 (3:05 EAST: OSE/EATCH/SMO: // LAB: END: // : NORTH: OCWN:
IALYSIS REQUESTED: INTAINER PRESERVATIVE MGP NAME .ASS ICED SM METALS .ASS ICED SS SEMIVOLATILES .ASS NONE STO9 CYANIDE .ASS ICED SM34 MERCURY BY COLD VAP delete
South end of old mound street plant in drive way 0-2 ft depth
Split Collected
BØ1 * Power Plant

AMPLE COLLECTED EY: Wes MEPal, John Peck

AFT U. Environmei	.S. ENVIRONMENT NTAL SERVICES O	FIELD SHEET AL PROTECTION IV. 25 FUNST	AGENCY/ REG On RD. Kansas	GION VII S CITY, KS ó	6115
: 91 4CTNG	: DSX44 SAMNG:	102 GCC: _ ME	DIA: SOIL	L: S P F D	********
	: LACLEDE COAL	GAS MO PROJEC	T NUM: A34	REF LATITUDE PT: LONGITUD	E:
	. LOUIS	OM LA8:	0/ BEG: 03// END:/	TE TIME 21/91 <u>/4:22</u>	FROM REF PT EAST: NORTH:
ALYSIS REG NTAINER ASS ASS ASS MMENTS:		STO9 · CY	TALS MIVOLATILES ANIDE RCURY	BY CULD VAP	delete.
	net of	2 old		street	!
	3¢1 ×	BØ2 * Powe	Mound Street - Plunt		,
•	Split C	ollected	·		

MPLE COLLECTED EY: Wea McCall, John Pock Petty Reports

FIELD SHEET U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 : 91 ACTNO: DSX44 SAMNC: 103 QCC: _ MEDIA: SDIL PL: S P F D MO PROJECT NUM: A34 PT: LONGITUDE: DATE OF TIME FROM REF PT

SE/BATCH/SMO: // LAB: ENU: 3/04/9/ : NORTH: DRET/SAROAD NO: ___ ALYSIS REQUESTED: PRESERVATIVE NTAINER MGP NAME 122 SM SS .SEMIVOLATILES 122 NONE STO9 CYANIDE ASS SY COLD VAP delete ICED 455 MENTS: JUFARE WILL COCKIED AT THE EAST CENTICAL PARCE OF THE TAME FARM - SEE MAP FOR LOCATION JAMPLE CENSISTS OF 1 ACIGICAT AT 1FT AMP I ALICLET AT 2 =T SPLIT SAMLE LAS COLLECTED

MPLE COLLECTED BY :

ENVIRONM		NTAL PROTECTION DIV. 25 FUNST	·		6115
: 91 ACTN	NO: DSX44 SAMNO	: 104 GC: _ ME	DIA: SOIL P	L: S P F D	
TIVITY CE	S: LAGLEDE CO	NL GAS MG PROJEC	R NUM: 234 P	EF LATITUD T: LONGITU	E:
CATION: S	ST. LOUIS /SMO:/_/ DAD NO:	MO LAB:	DA BEG: 03/2 END:/_	TE TIME 191 <u>/7:25</u> _':_	FROM REF PT EAST: NORTH:
ASS		TIVE MGP NA SM ME SS SE STO9 CY SM34 ME	TALS MIVOLATILES	BY-COLD-VA	-cielotto
MMENTS:	Boring	the two tonk to Location collected.	•		north

		•
1FT	II C THUTBONNENTAL	FIELD SHEET
		L PROTECTION AGENCY, REGION VII V. 25 FUNSTON RD. KANSAS CITY, KS 66115
91 ACTN	IO: DSX44 SAMNO: 10	O5 CCC: _ MEDIA: SDIL PL: S P F D
:MCITA:		MO PROJECT NUM: A34 PT: LONGITUDE:
MPLE DES:	B07	MO BEG: 03/201/91 /2:00 EAST: LAB: END: 1051 : NORTH: DCWN: 2-2 for
CATION: S	T. LOUIS	MO 6EG: 03/81/91 /3:10 EAST:
SE/BATCH/	'SMO:/_/_	LAB: END: PSI : NORTH:
ORET/SARD	DAD NO:	DOWN: 2-7
ALYSIS RE		•
	PRESERVATIVE	MGP NAME
ASS	ICED	SH METALS
ASS	ICED	SM METALS SS SEMIVOLATILES
ASS	NONE	STO9 · CYANIDE
ASS	ICED	STO9 CYANIDE SH34 MERCURY BY CULD VAP clickete
MMENTS:		
	Between.	tank 7-8
	, -	
	le bou	ing & safety logitard #1

and the control of the control of the control of the control of the control of the control of the control of t

171

	ENVIRONMENTAL L SERVICES DIV.				66115
': 91 ACTNO: D	SX44 SAMMO: 10	S GCC: _ MED1	A: SOIL	PL: S P F ()
CATION:	ACLEDE COAL GAS	MO PROJECT			JDE:
.MPLE DES: _/3 CATION: ST. L .SE/BATCH/SMG: 'ORET/SAROAD N	OUIS 0:	MO LAS:	### BEG: 03/ ENO:/	DATE TIME 181/91 <u>/2</u> : <u>/</u> 1	FROM REF PT PEAST: NORTH: DOWN:
IALYSIS REQUES	TED: PRESERVATIVE ICED	MGP NAMI SM META SS SEM	E ALS IVOLATILES NIDE		ar clileti
IMMENTS:	brung	* safet) log	book #	L

South of little tankform near

AMPLE COLLECTED BY : FRIE (1) (4)

	S. ENVIRONMENT TAL SERVICES D				115
: 91 ACTNO:	DSX44 SAMNE:	107 GCC: _ ME	DIA: SOIL	PL: S P F D	
SATION:	LACLEDE COAL Y <u>louic</u>	MO PROJEC			
APLE DES: CATION: ST. SE/BATCH/SM DRET/SAROAO	S/6 Louis 0:/_/ No:	MO LAB:	8EG: 03 END:	DATE TIME = 101/91 /5:/5 = N	POM REF PT AST: ORTH:
ALYSIS REQU VTAINER ASS ASS ASS	PRESERVATIV ICED ICED NONE ICED	SM 'ME	TALS Mivolatiles	BY COLD VAP	delete.
	20ft. Lu boa				

		PROTECTION AGENCY, REGION VII 25 FUNSTON RD. KANSAS CITY, KS 66115
: 91 ACTNO: 05	X44 SAMNG: 108	CCC: _ MEDIA: SOIL PL: S P F D
	ACLEDE COAL GAS	REF LATITUDE:
MPLE DES:	/_/	MO BEG: 03/01/91/5:45 EAST: LAB: END: 104/ : NORTH: DOWN: 0-2/1
ALYSIS REQUEST		
	PRESERVATIVE	MSP NAME
ASS	ICED	SM \ METALS SS \ SENIVOLATILES
ASS	NONE	STOO CYANTE
ASS	ICED	SH34 HERCURY BY COLD VAP () Lete?
MMENTS:	5' north	of little tankform.

FIELD SHEET

IMPLE COLLECTED BY : 12/1/ 1974 (411

AFT		FIELD SHEET			
	J.S. ENVIRONMENTAL Ental Services Div				6115
91 ACTNO	D: DSX44 SAMNO: 10	9 QCC: _ MED	IA: SDIL	PL: 5 P F D	
:NCITA:	S: LACLEDE COAL GA	MO PROJECT	NUM: A34		DE:
4PLE DES: CATION: S' SE/BATCH/S CRET/SARO	T. LOUIS SMO:/_/ AD NO:	MO LAE:	9EG: 03 END:	DATE TIME 191191 <u>/6:57</u> 1061:_	FROM REF PT EAST: NORTH:
ALYSIS REF VTAINER ASS ASS ASS ASS	QUESTED: PRESERVATIVE ICED ICED NONE ICED	MGP NAM SM MET SS SEM STO9 CYA SH34 MER	ALS IVOLATILES NIDE	·	-delati
4MENTS:	110' m	ith of.	littl	ta, ky	Lauin
	su lo	gbook >	42	-	

ENVIRGNMEN	S. ENVIRONMENTA TAL SERVICES DI	V. 25 FUI	ION AGENCY, NSTON RD. KAI	NSAS CITY,	KS 66115	
r: 91 ACTNE:	DSX44 SAMNO: 1	10 QCC: _				
TIVITY CES:	LACLEDE COAL G	AS		REF LAT 4 PT: LON	ITUDE:	
	#/7 LOUIS 10:/					
NALYSIS REQU INTAINER LASS		MGP SM SS	NAME METALS SEMIVOLATIL			,
OMMENTS:	S of si	on r	achord	fy trace		

Brown / Haxes SH

AMPLE COLLECTED BY : _

RAFT

CRAFT

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

ENATYONNEN	INE SERVICES DIV	. 23 FU	usion ko-	VWUJWJ	CTIIN K2	00113	
FY: 91 ACTNO:	DSX44 SAMNO: 11	1 900: _	MEDIA: SO	IL P	L: S P F D		
	LACLEDE COAL GA		JECT NUM:		EF LATITUD T: LONGITU		
;AMPLE DES: _	RDD			7. Ø DA	TE TIME	FROM REF PT	
		MO	BE	s: 03/P	1791 ///2. 7	FEAST:	
CASE/BATCH/SM	. LOUIS	_ LAB:	ENG	D:/#	21_ 27:30	NORTH:	
TORET/SARDAD	NO:	_		E	SA	NORTH:	pt-
ANALYSIS REQU	JESTED:						
CONTAINER	PRESERVATIVE	MGP	NAME			·	
JLASS	ICED	SM	METALS				
GLASS	ICED	S \$	SEMIYOLA'	TILES			
LASS	NONE	ST09	CYANIDE		alele de		
JLASS	ICED	'SM34	MERCURY-		BY COLD VI	TP	

:OMMENTS: .

S. of mullingly 180' E. of R.R. truck laybook #2

Brown Hayes SH

TAMPLE COLLECTED BY : TPC

TAAT

PIELD SHEET

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ΞN	VIF	RON														ST											6 :	611	; 			
Y: 3	1 .	CT	NC:		25)	44	5	ΔM	NC	:	112	<u> </u>	CC	:	_	ME	DIA	\:	SB	IL		9 [.:	S	ρ	F	٥					
CTIV										L ==				2 6	เขา	EC	T N	IUM	:	434	4						Dē	•				
LHPL																				ス	څ.	DA1	ΓE		ŢI	ME		= R O	4 2	EF	PT	
DCAT												i40						3	ΞG	: (33.	101	119	71	<u>/~</u>		10	EAS.	Γ:			
ASEI	AE.	TCH	13	10	: _			/_	/			_	L	AE	3:		_	Ξ	ND	:		ني	71 _		47		7 0	NOR	TH:			
TORE	T/:	5 4 R	DAI) ·	10	: -			-								_					13	¥	,	4	14	_	NOR'	Y :	0		4
NALI	SI	ऽ २	EQI	JE!	STE	٥:																									•	
THE	INI	ER				5 5 E	SE	RV	A T	۲v	E		MG	P		MA	ME															
LASS	5				•	CE	D						SM	}		ME	TAL	. S														
LASS	5					CE	D						SS			SE	MIN	OL	AT	IL	E S											
LASS	5				ł	NON	ΙE						ST	09	•	CY	AN]	EDE														
LASS	5					I C E	D					s	SM	34	-	ME				\ll	=	ر <u>ح</u>	BY	Ç	JLI) y	AP					

:STMBMES:

for w. of with the

logbook #2

Borr Low 12

Hayes Brown SH

TAMPLE COLLECTED BY :

	FIELD SHEET ENVIRONMENTAL PROTECTION AGENCY/ REGION VII L SERVICES DIV. 25 FUNSTON RD. KANSAS CITY/ KS 66115	
	SX44 SAMNO: 231 GCC: _ MEDIA: WATER PL: S P F 0	
ACTIVITY DES: L	ACLEDE COAL GAS REF LATITUDE: 1D PROJECT NUM: A34 PT: LONGITUDE:	
LOCATION: ST. L LOCATION: ST. L CASE/BATCH/SMO: TORET/SARDAD N	ONTE FROM REF PT OUIS MO BEG: 03/04/9 : EAST:	
	PRESERVATIVE MGP NAME ICED WV VOLATILES	
COMMENTS:	Coll sample -ottertes sample at 36 test. coation = 64 test 1 from certer	
	STATIC WATER LEVEL: 24 FT	<i>K</i> S
	DEPTH OF WELL: 26 FT.	
	= 7.58 $5 = 560$	
1/1/9	6 –	

	PIE - ENVIRONMENTAL PA AL SERVICES DIV.		AGENCY, R		56115
TY: 71 ACTNO:	DSX44 SAMNO: 202 G	CC: _ M5	DIA: WATER	PL: S P F D	
ACTIVITY DES:	LACLEDE COAL GAS) PROJEC	T NUM: A34	REF LATITUDE PT: LONGITUE	DE:
IAMPLE DES: LOCATION: ST. DASE/BATCH/SMG TORET/SAROAD	LOUIS MO	LAB:	8EG: 03 END:	3.06/9/ (FE OATE TIME 1.01/91	FROM REF PT EAST: NORTH: DOWN: 33
2 VOA VIALS CUBI ;LASS , GZ PLASTIC CUBI	PRESERVATIVE ICED 5 ML HNG3 ICED FILTER, HNG3	WW MI WS SI WO7 II	EMIVOLATILES H dissolved	METALS AL SY COLD VA	edelete-
	Coll' San Tennival F BR Frack Center or Depth to Depth of	RAS	soc. / 78 Ft. Nu/land	South	of the
- em)	7 = 7.8/1 = 7°C				

Williams

CRAFT		IELD SHEET		
	S. ENVIRONMENTAL			
ENVIRONMEN	TAL SERVICES DIV.	25 FUNSTON RD	. KANSAS CITY,	KS 66115
=Y: 91 ACTNO:	DSX44 SAMNC: 203	QCC: _ MEDIA:	WATER PL: S P	F D
	LACLEDE COAL GAS		REF LAT	
LOCATION: _C	t. Louis	40 PROJECT NUM	1: A34 PT: LOM	GITUDE:
SAMPLE DES: _			03/47/4/92	IME FROM REF PT
CATION: ST.	LOUIS	6 OM	EG: 63/01/91	
CASE/BATCH/SH	IC:/_/	LAB: E	ND: / / 7	V: 30NORTH:
STORET/SAROAD	NO:			DOWN: 473"
ANALYSIS REQU	IESTED:			
	PRESERVATIVE	MGP NAME		
2 VOA VIALS	ICED	WV .VOLATIL	.ES	
CUBI	5 ML HNO3	WM METALS		
FLASS	ICED	WS · SEMIVOL		
	FILTER, HNO3	WO7 -IH DISS		
CUBI	NAOH	HT09 CYANIDE	TOTAL	inete.
CUBI	5 ML HNO3	WM34 MERCURA	SY COL	D VAP de le te.
COMMENTS:	ARMY CORPS O	F ENGINEERS	•	
COMMENTS:		11,		
	Engineering	lest w	10//	
	Engineering Location =	147 fee	of north	of the
	center	of Mu	Man aher	Start and Sourcete floor
			wan phy	suckt and
	40-10	ef wes,	1 0/0	oncrete flias
	leve	<i>/.</i>	- ,	
	Depth de		72"	
	Depra of	WEN = 4,	<i>. 3</i>	

· p#:6.32. (ano. - 1500 TEMP . - 9°C.

AMPLE COLLECTED BY: Robert / Williams

Static water lavel = 27'.3"

;RAFT	=	FIELD SHEET	
		PROTECTION AGENCY, REGION VII	
		25 FUNSTON RD. KANSAS CITY, KS 65115	
=Y: 91 ACTHG: 3	SX44 SAMNG: ?74	4 RCC: _ MEDIA: WATER PL: S P F D	
CTIVITY DES: L	_		
TXT : NEITABL	Louis	MB PROJECT NUM: A34 PT: LONGITUDE:	
	a	JATZ GENE FROM REF PT	
JAMPLE DES: LOCATION: ST. L	OUTS	NO REC. DIAGO FACT.	
145F/8ATCH/SMG:	/ /	MO 3EG: 03/01/91 : EAST:	
CAOSES/TEROTE	0:	DOWN: 48	
INALYSIS REQUES	TED:	·	
-	PRESERVATIVE	MGP NAME	
I VOA VIALS		WV / VOLATILES	
cuai	5 ML HNO3	WM METALS	
JLASS	ICED	WS SEMIVOLATILES	
	FILTER, HNO3	WO7 -IH DISSOLVED METALS	
	NAOH 5 Ma ung 7	WT09 CYANIDE, TOTAL BY COLD VAP Lilitta	
CUBI	5 ML HNO3	WH34 MERCURY BY CULD VAP COLOCIO	
COMMENTS:	10 64.004		
	• •	of Engineers	
	Facines 1	ing Trest well	
	Digine	ing IRST Well	
	11. 1:1		
	_ocation	= 216 feet north of well	
	^		
	(Sample	#203 was collected fearing	
	3625	et north of the center of	
_			
	Mulland	thy street and 25 feet we	
	2 /2	ing shout and as feet we	2
	1- X10	Courseta Elas	
	Depth o	of well= USI	
	Shi	of well= 481	
, /	(0 /	- CAG	
. 247	- 6.21.		
	000.		
: /EMP =	7		
Conol. =	400 um ho	25	
A IPLS COLLECT	17 14 : <u>Z</u> o	operts/Williams	

٠	2	•	=	T
	er.	-	_	1

	S. ENVIRONMENTAL TAL SERVICES DIV				56115
TY: 31 ACTNO:	DSX44 SAMNE: 20	5 QCC: ME	DIA: WATER	PL: S P F D	
ACTIVITY DES:	LACLEDE COAL 3A	S 40 PROJECT	T NUM: 434	REF LATITUD: PT: LONGITU	E:
SAMPLE DES: _A LOCATION: ST. CASE/BATCH/SMO STORET/SAROAD	wouls 1. / / / / / / / / / / / / / / / / / / /	17e 40 LAB:	9EG: 03 =NO:	ORTE TIME 194191 01:30	FROM REF PT FAST: NCRTH:
ANALYSIS PEQUATONIAINER 2 VOA VIALS CUBI GLASS 4 DI PLASTIC CUBI CUBI	PRESERVATIVE	WW . VOI . WW . ME' . WS . SEI	LATILES TALS Mivolatiles - Dissglved Anide, Tota	HETALS "	politica e
Sampo	linsate of leo-usi	l auger	tean o	los/levs	l groon decon

CRAFT

FIELD SHEET

	1.5. ENV					
ENVIRONME						
. 14 46716						

Y: 31 ACTNO: DSX44 SAMNO: 206 QCC:	_	PL: S P F D	
CTIVITY CES: LACLEDE COAL GAS GCATION: 57. Louis MO 6	PROJECT NUM: A34	REF LATITUDE:	
AMPLE CES: _OCATION: ST. LOUIS MO _ASE/BATCH/SMG:// L/ TORET/SAROAD NO:		DATE JIME FROM 3/01/91 : EAST 19/9/10:27 NORT	REF PT

NAME

INALYSIS REQUESTED:

SNTAINER PRESERVATIVE MGP VOA VIALS ICED WV

ICED WY VOLATILES
5 ML HNO3 WM /METALS

LASS ICED WS /SEMIVOLATILES
OZ PLASTIC FILTER, HNO3 WO7 , IH DISSOLVED METALS

OZ PLASTIC FILTER, HNO3 WO7 , IH DISSOLVED METALS
UBI NAOH WT09 CYANIDE, TOTAL

CUBI NAOH WTO9 CYANIDE, TOTAL
CUBI 5 ML HNO3 WH34 MERCURY BY COLD VAP (1) (1)

COMMENTS:

CUBI

COLLECTED FROM ARMY CORPS OF ENGINEERS

LOCATION: 20 FT. WEST OF CONCRETE
LEVEE WALL AND ~ 21/8 FT

N. OF THE CENTER OF

MULLANIAY 57.

DEPTH OF WELL: 52 FT STATIC WATER LEVEL: R9 FT.

. PH = 6.14. , CCNO. = 1600 инфозі . ТЕПР. = 11°С

Le Michael Manning

TAMPLE COLLECTED BY : ROBERTS & WILLIAMS

The let

	• • • •		•				
ENVIRONMENTA	ENVIRONMENTAL L SERVICES DIV.	PROTECTI 25 FUN	ON AGENCY			6115	
FY: 91 ACTNO: 1	35X44 SAMNO: 20	gcc:	MEDIA: WA	TER PL:	SPFD		
CTIVITY DES: I	ACLEDE COAL GAS	s S		REF	LATITUDE		
AMPLE DES: _CCATION: ST. CASE/BATCH/SMO _TORET/SAROAD	LOUIS:	MO LAB:	B E G END	DATE: 03/01/ : 3/9/	A 10:22	FROM REF EAST: NORTH: _ DOWN: _	PT <u>S2</u> FT
VOA VIALS LUBI GLASS OZ PLASTIC	PRESERVATIVE ICED 5 ML HND3 ICED FILTER, HND3	' WV ' ' WM ' ' WS ' ' WO7 ' '	VOLATILES METALS SEMIVOLAT TH DISSOL	ILES VED META		Olele	ete.

COMMENTS:

COLLEGED FROM ARMY CORPS OF ENSINERS

LOCATION: 20 FT. WEST OF CONCRETE

LEVEE WALL AND ~ 2118 FT

NORTH OF THE CENTER

OF MULLANPHY ST.

DEDTH OF WELL: 52 FT

SFATIC WATER LEVEL = 29 FT.

PH = 6.14'

COND, = 1600 Unhos.

TENP. = 11'C'

SAMPLE COLLECTED BY : ROBERTS + WILLIAMS

Their G. H.

RAFT

FIELD CHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 46115 FY: 31 ACTNO: ISX44 SAMNC: 207 ACC: MEDIA: WATER PL: S P F D CTIVITY DES: LACLEDE COAL GAS LECATION: SK Lanie 40 PROJECT NUM: 434 PT: LCNGITUDE: AMPLE DES: DATE TIME FROM REF PT BEG: 03/01/94 LOCATION: ST. LOUIS MO :TZA3 #SE/BATCH/SMC: ____/_/__ LA3: ____ END: 03/09/2/ TORET/SARDAD NO: ____ INALYSIS REQUESTED: ONTAINER PRESERVATIVE HGP NAME _ VOA VIALS 1 MA **/ VOLATILES** CUBI 5 ML HNO3 · METALS · WW LASS ICED · WS , SEMIVOLATILES. OZ PLASTIC FILTER, HNG3 , ¥07 'IH DISSOLVED METALS SWT09 'CYANIDE, TOTAL CUBI NAOH CUBI 5 ML HNO3 ? Oozglass Icho COMMENTS:

Fiel D Blank

TAMPLE COLLECTED BY: Williams/ Roberts

_RAFT	F	IELD SH	EET	• •			
PNVIKUNMENT.	ENVIRONMENTAL AL SERVICES DIV.	25 511	N.C.TON	00 W 44			66115
FY: 91 ACTNO:	DSX44 SAMNO: 207	9cc: 1	MEDIA	: WATER	PL: S	PFD	
LOCATION:	LACLEDE COAL GAS	; Mo pro	JECT N	UM: A34	REF L	ATITUD. UTIDKO.	E:
AMPLE DES: CATION: ST. CASE/BATCH/SMO TORET/SAROAD	LOUIS	MO LAS:		BEG: 0 END: _	75/05/9/ 94 TE 3701791 _//	PEIME ZZ:00	FROM REF PT EAST: PNORTH: DOWN:
GLASS OZ PLASTIC	PRESERVATIVE ICED 5 ML HND3 ICED FILTER, HND3	. WY . WM . WS . WO7 . WID9	VOLAT METAL SEMIV 'IH DI CYANT	OLATILE SSOLVED De Tot	METALS		edelli.
COMMENTS:	GEO probe Poured prod COLLECTED BY					es .	and

PH = 1.00.

TAMP = 4°C.

Cono. = 16 ampos

SAMPLE COLLECTED BY: Boser's Williams

LRAFT		FIELD SH	-		
	U.S. ENVIRONMENTAL BENTAL SERVICES DIV	. 25_EUI	NSTON RD. KA		
FY: 91 ACTN	D: DSX44 SAMNO: 20		MEDIA: WATE	R PL: S	P F D
CTIVITY DE	S: LACLEDE COAL GA	S MO PRO	JECT NUM: A3		ATITUDE:
AMPLE DES: _CCATION: S CASE/BATCH/ TORET/SARC	ST. LOUIS SMO:/_	MO LAS:	BEG: END:	DATE 93/01/91 93/01/91	TIME FROM REF PT EAST: NORTH: 18:30 DOWN:
VOA VIALS	PRESERVATIVE SICED 5 ML HNO3 TLED	MGP YWY WH WS	NAME VOLATILES Metals Semivolatil		1
LOZ PEAST	IC FILTER/HNUS	¥07	TH DISSULVE		V

COMMENTS:

Trip Blank

SAMPLE COLLECTED BY: Williams Riberts

FIELD SHEET U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV." 25 FUNSTON RD. KANSAS CITY, KS 66115 FY: 91 ACTNO: DSX44 SAMNC: 301 GCC: _ MEDIA: WATER CTIVITY DES: LACLEDE COAL GAS PROJECT NUM: A34 FROM REF PT LOCATION: ST. LOUIS CASE/BATCH/SMO: ____/_/_ STORET/SAROAD NO: ANALYSIS REQUESTED: CONTAINER PRESERVATIVE MGP NAME 2 VOA VIALS 5 HL HNO3 CUBI

WS

CUBI 5 ML HNO3

3 40 ML VOA VL3 ICEB

8002 9/455 PE

FILTER, HNO3

COMMENTS:

4 OZ PLASTIC

SLASS

CUBI

SW sample

Abandoned Pump lhouse, Collected

At the SE Corner of pump

house

HYDROCARBONS, TOTAL PETROL,

pH = 8.15 Temp. = 7°C Cord. = 440 amhos

SAMPLE COLLECTED BY: Deberts Williams

5 4 F T

FIELD SHEET

L.S. ENVIRONMENTAL PROTECTION AGENCY/ PEGION VII EMVIFORMENTAL CERVICES DIV. 15 FUNSTER RD. KANSAS CITY, KS 66115 CTIVITY DES: LAGLEDE SCAL GAS ECCATION: ST. LOUIS MO BEG: 63/01/91 END: _/_/_ 1/2:43 152/3ATCH/SMC: ____/_/__ LAB: __ T026T/54R040 %0: -NALYSIS REQUESTED: PRESERVATIVE ONTAINER AGP. NAME VK VOA VIALS ICED **VOLATILES** V 5 ML HNO3 · MM METALS / LASS ICED SEMIVOLATILES / . WS DZ PLASTIC FILTER, HND3 IH DISSOLVED METALS **M07** CUBI NACH CYANIDE, TOTAL TUBI 5 ML HNO3

OMMENTS:

80 02 9/ess (DZ

Duplicate sample of \$\$ 30/

HYDROCARBONS, TOTAL PETROL

Joans Williams

imple Coulisted by :

TEELS SHEET U.S. ENVIRONMENTAL PROTECTION AGENCY/ REGION VII TY: 91 ACTNO: DSX44 S4MNO: 302 QCC: _ MEBIA: WATER CTIVITY SES: LACLEDE COAL JAS AMPLE DES: LOCATION: ST. LOUIS 10 1452/3ATCH/SMO: ____/_/__ END: __/_/_ /2:/5 NCRTH: TORET/SARGAD NO: :NALYSIS REQUESTED: CNTAINER PRESERVATIVE MGP VOA VIALS **VOLATILES** * 5 ML HNO3 LASS SEMIVOLATILES / W07 OZ PLASTIC IH DISSOLVED METALS FILTER, HNO3 CUBI CUBI HYDROCARBONS, TOTAL PETROL DMMENTS:

SW sample

Collected about MOFF upsteram

of NE Corner of the abandon of

pump house.

1/2 7.87. 1/emp. = 6° C 310, = 450 cm/105 PLE SOLLESTED IN: Boperts Williams

FIELD SHEET U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 -CTIVITY DES: LACLEDE COAL GAS REF LATITUDE: AMPLE CES: LOCATION: ST. LOUIS END: __/_/_ /4: @ NORTH: ASE/SATCH/SMG: ____/__ LAB: ___ TORET/SARGAD NO: INALYSIS REQUESTED: PRESERVATIVE CHTAINER MGP NAME _ VOA VIALS · VOLATILES 5 ML HN03 CUBI LASS 'SEMIVOLATILES WQ7 OZ PLASTIC FILTER, HNO3 ·IH DISSOLVED METALS CUBI WT09 -CYANIDE, TOTAL WH34 HERCURY BY COLO VAPOLETE 5 ML HN03 CUBI W486 - HYDROCARBONS, TOTAL PETROL 8002 9/495 **OMMENTS:**

SW SAMPLE Miss. River

Collected 330 feet South of the southeast corner of the abandured pump house

1/270

TRAD = 5°C'

mad = 360 mm 205

TAMPLE COLLECTED BY: Hoberts Brawn

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 TY: 91 ACTNO: DSX44 SAMNO: 304 QCC: _ MEDIA: WATER PL: S P F D MO PROJECT NUM: A34 PT: LONGITUDE: SAMPLE DES: TIME BEG: 93/01/91 LOCATION: ST. LOUIS MD EAST: CASE/BATCH/SMO: ____/__ LAB: ___ END: 7/7/9/8:55 NORTH: STORET/SAROAD NO: __ INALYSIS REQUESTED: CONTAINER PRESERVATIVE MGP NAME 2 VOA VIALS ICED WV VOLATILES V CUBI 5 ML HNO3 WW METALS V JLASS ... ICED WS SEMIVOLATILES / 4 OZ PLASTIC FILTER, HNO3 W07 IH DISSOLVED METALS CUBI NAOH WTO9

HYDROCARBONS, TOTAL PETROL /

COMMENTS:

CUBI

SURFACE CATER SAMPLE FROM THE

ILLINOIS AMERICAN PRINKING WATER

INTAKE IN E. ST. COUS.

, PH . MESTON TY THE LEMNES.

TEPP. 3°C:

5 ML HN03

FIELD SHEET

· F				7	311551					
	U-3-	ENVIRGNMEN	ITAL	PROTE	RCTION	AGENCY	P REGI	ION VII		
EMVIRGNM	ENTAL	SERVICES	DIV.	25	PUNSTO	N RO.	KANSAS	CITY,	KS	56115

=y: →1 ACTNG: @SX4	4 SAMNC: CC: _	MEDIA: WATER PL: S P F D	
.CTIVITY DES: LACL		REF LATITUDE JECT NUM: A34 PT: LONGITUD	
JAMPLE DES: LDCATION: ST. LOUI DASE/BATCH/SMC: TORET/SARDAD NO:	S MO LAB:	DATE TIME BEG: 03/01/91 END @ / ^/ (1/ -: 55	FROM REF PT EAST: NCRTH:
TLASS IC TOZ PLASTIC FI CUBI NA	ESERVATIVE MGP ED WW ED WS LTER, HNO3 WO7 BH WT09 ML HNO3 W86	METALS/ SEMIVOLATILES / IH DISSOLVED METALS / CYANIDE, TOTAL / HERGURY BY COLD VAP	-delete pe

:DMMENTS:

OUPLICATE SURFACERATER SATTPLE FROM THE LLINGS HITERICAN, PRINKING -LATER MITAKE IN E. T. LOCIS

TEMPO 3°CI

JAMPLE COLLECTED BY :

,

13 45 7	<u>چ</u>	FIELD SHEET
ENVIR	The state of the s	PROTECTION AGENCY, FEGION VII , 25 FUNSTON 30. KANSAS CITY, KS 66115
Y: 31 A	TNO: 25x44 S4MNC: 401	CC: _ MEDIA: SOIL PL: S P F D
CTIVITY SCATION	CES: LACLEDE COAL GAS	REF LATITUDE:
LOCATION LASE/BAT	ES: : ST. LOUIS Ch/SMO:/_/ ARDAD NO:	MO 3EG: 93/01/91 : EAST: LA3: END: _/_/ 75:30 NORTH: DOWN: 0-6
ONTAINE LASS GLASS -LASS LASS	NOME ICED ICED Nome	SV - VOLATILES SM - METALS SS - SEMIVOLATILES STO9 - CYANIDE
SLASS	ICED	SH34 HERCURY BY COLD VAP COLD VAP

COMMENTS:

SKOIMENT SAMPLE / GAB Collected about 150ff downstream from the SE corner of the abondoned pump house.

SV54 'HYDROCARBONS, TOTAL PETROL

SIET

FIELD SHEET

ENVIRON	U.S. ENVIRO Mental Servi	CES DIV. 2	5 FUNST	=	PEGION VII SAS CITY, KS	ó ó115
-Y: 71 ACT	NO: DSX44 SA	MNG: 474 20	C: ME	DIA: SOIL	PL: S P F D	
	St. Louis		PROJEC	T NUM: A34	REF LATITUD PT: LONGITU	
SAMPLE DES	:			a	FORE/PEINE	FROM REF PT
	STUCL .T?	MO		3EG: -0	3/01/9 1;	EAST:
TASE/EATCH	// SMG:/	_/	LAB:	END: _	3/01/91 _//_ <i>15_3</i> 2	NORTH:
	יפר פגני:					SOWN: O.6
ANALYSIS R	EQUESTED:					
CONTAINER	PRESER	VATIVE N	IGP NA	ME		
ILASS	NONE	· s	v vo	LATILES		
GLASS	ICED	7 S	M 'ME	TALS	•	
GLASS	ICED	, 5	S 'SE	HIVOLATILE	S	
LASS	NONE	/5	T09 -C1	ANIDE		1 1 1
GLASS	ICED		M34 HE	RCURY	BY COLD VI	recelet
GLASS	ICED	45	1754 . HY	DROCARBONS	. TOTAL PETRO	IL .

COMMENTS:

Duplicate of #401

TAMOLE COLLECTED BY: To berts/Williams

TALET

FIELD SHEET

U.S. ENVIRONMENTAL PROTECTION AGENCY, REGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115

FY: 91 ACTNO:	: DSX44 SAMNO: 40	2 1CC: _	HEDIA: SOIL	PL: 3 9 =	j)	
230 YTIVITO.	LACLEDE COAL GA		JECT NUM: A3	REF LATITU	~ ~ ~~ ~~	
:AMPLE DES:			(03/06/9/ (PC) HE	FROM REF PT	
LOCATION: ST.	LOUIS	หอ	BEG:	03/01/91 :		
CASE/BATCH/SR	1G:/_/_	LAB:		///Z:3		_
.TGRET/SARDAD	. NO:	_		/-	25MN: 0-6	,
:NALYSIS REQU	JESTED:					
CONTAINER	PRESERVATIVE	НGР	NAME			
LASS	NONE	SV	"VOLATILES		•	
JLASS	ICED	SM	METALS			
TLASS	ICED	SS	SEMIVOLATIL	.ES		
				- -		

COMMENTS:

SEDIMENT SAMPLE/GRAB

(allectro about 60 fret south of

SW Samp & # 302.

SW34 MERCURY BY COLD VAP delete "
SV54 HYDROCARBONS, TOTAL PETROL

Brown Williams

FIELD SHEET

		PROTECTION AGENCY, 25 FUNSTON RO. KAN	
-Y: /1 -CTNG:	35X44 SAMNO: 403	OCC: _ MEDIA: SOIL	PL: S P F D
.ccation:	LACLEDE COAL GAS	TO PROJECT NUM: 434	REF LATITUDE:
JAMPLE JES: _ LOCATION: ST. JASE/BATCH/SM JTORET/SARDAD	LOUIS 0:/_/	MO 3EG: 0 LAB: END: _	07,42 PIME FROM REF PT 3401791 : EAST: 1 /4: /2 NORTH: 20WN: 0-6
ILASS GLASS	PRESERVATIVE NGNE ICED ICED	SV VOLATILES SM METALS SS SEMIVOLATILE	ES BY COLD VAR ChileTee
COMMENTS:	SEDIM	KNT SAMPLIE	- Carab Miss. River
_	of the s	330 fre f so southerst com pump house	with (downsterm) ver of the

CHAPLE COLLECTED IN: Schools Brown

FIELD SHEET J.S. ENVIRONMENTAL PROTECTION AGENCY, PEGION VII ENVIRONMENTAL SERVICES DIV. 25 FUNSTON RD. KANSAS CITY, KS 66115 🛴 AMPLE DES: 80,000 gol Tank #80-1 BEG: 0344791 7:45 EAST: END: _/_____ NORTH: :ASE/BATCH/SMO: ____/_/___ LAB: ___ TOPET/SARDAD NO: ANALYSIS REQUESTED: CHTAINER PRESERVATIVE HGP MAME , UV **'VOLATILES SEMIVOLATILES** 40 ML VOA VLS ICED -WYGG -HYDROCARBONS, TOTAL PETROL.

COMMENTS:

Sample collected out of side value of an 29,000 gal. tank # 80-1. It is located on the S.W. corner of the tank farm.

His Constitution sample - Oil Simple

INDLE COLLECTED ET : Klieb A. Brown

DRAFT

中间压力 与相称的不

TO SEE ENVIRONMENTAL PROTECTION AGENCY. REGION VIL PAYIRONMENTAL SERVICES DIV. 195 FUNCTION RD. KANSAS CITY, KS 85113

TY: MI SUINU: DSX44 SSAMO: THE OUT: P MEDIA: WAIER MY: S P F D

OFFICED BY SECTION FOR THAT YOUR

BEE 工商1141108年

. SAMPLE DEST LACLEDE COME GAS SUITEDE SAMPLE DATE TIME CROM REF DI

COUNTRIES AND MA 860: 03/08/21 _ : FAST: LAB: FMD: / / ... NORTH:

STORET/SARUAD NO: []

DOMN:

AMALYSIS REQUESTED:

CONTAINER

PRESERVATION

MUP MAME

2 YOA VIALS

TCED TUED

WV VOLATILES

GLASS CUBI

5 ML HN03

98 SEMICYDUATIOLES

4 DZ PLASTIC

HM METALS

FILTER HND3

TH DISSOLVED BETALS **907**

CHRI

NAGH

W109 CYANDE: TOTAL

CHRI

5 ML HN03

COMMEN(5:

CHAIN OF CUSTODY RECORD ENVIRONMENTAL PROTECTION AGENCY REGION VII

ACTIVITY LEADER(P	rint)	,	NAME	OF SURVEY	OR ACTIVITY		-	1		DATE OF COLLECTION SHEET DAY MONTH YEAR OI /		
	CONTENTS OF SHIPMENT						Carl Cos S. TE. DAYON MONTH YEAR					
SAMPLE	J-//GA		OF CONTAIN		VOA SET	S	AMPI	LED M				
NUMBER	<u> </u>	SOTTLE SERS OF CONT	BOTTLE AINERS PER SA	30TTLE	(2 VIALS EA)	water	Š	sediment	Š	condition of samples upon receipt. Other sample numbers etc.)		
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DESCRIPTION OF S	SHIPMENT				MODE OF SH	IPM	ENI	1				
Ji33 PIECE(S),	CONSISTING ()F	SOX(ES)		сомм		AL (CARR	IER _			
Lice Ches	T(S). OTHER _				COURI		ากพ	VEVE	n	(SHIPPING DOCUMENT NUMBER)		
PERSONNEL CUST	ODY RECOR)	·		7					(Shirring Document Number)		
RELINGUISHED BY	Y (SAMPLER	DAT	E TIM		CEIVED BY			1 /	٠ ,	REASON FOR CHANGE OF CUSTODY		
Tutte	Bros		9 11		Jelle.	-44		يرس	u.	Manager TO EPA Lib		
SEALED STAND	UNSEAL	DAT	E TIM		CELYED BY		UI	NSEA	LED	REASON FOR CHANGE OF CUSTODY		
1 Willian	obele	3/8	1/11/09	2KX	1 cm		X	Tan	12	I mansport to EPA		
SEALED_	UNSEAL		/91		EALED BY		U	NSZ	LED	DEASON FOR CHANGE OF CUSTODY		
2-1	Fifty.			05-5	Jun 18	سند پیرم	m	اسا		Kica W. EPA Lake		
SEALED	UNSEA	LED /	1"		EALED	,	U	NSE	ALEC			

MAIN OF CUSTODY RECORD (ENVIRONMENTAL PROTECTION AGENCY REGION VII

ACTIVITY LEADER(PI	rint) Pontura		NAME CA N	:	OR ACTIVITY	,	<i>(a)</i>	٢	D	ATE OF COLLECTION 9/ DAY MONTH YEAR	SHEET /	
CONTENTS OF SHIPM				22252			, 0, 3			DAY MONTH YEAR	/11/	
			-	No. 64 S		ŞΑ		D MED	iA j	RECEIVING LABORATO		
SAMPLE !	CUSITAINER	122			. DA SET	1 5		QuSI	other	REMARKS OTHER INFORI	n receipt.	
	WIM66	RS OF	(, <u>), </u>	2011	- a			इ इ		other sample numbers, etc.)		
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ESCRIPTION OF S	HIPMENT				MODE OF SH	IIPMI	ENT			V		
PIECE(S) C	CHISISTING C	:			СОММ	ERCL	21 CA	RRIF	R			
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:CE CHEST	137 UTHER _				SAMPI	LER C	ONVE	EYED		SHIPPING DOCUMENT	NUMBERI	
ERSONNEL CUSTO	DDY RECORD)										
ELINQUISHED BY	(SAMPLER)	0/	ATE TH	प्रदेश विह	CEIVED BY		.3,	1141	i)	REASON FOR CHANGE	OF CUSTOD	
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ELINQUISHED BY			• · c.	. •	SE VED BY					REASON FOR CHANGE	OF CUSTOD	
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SEALED ELINQUISHED BY	- FSEAL		 		CEIVED BY		214	<i></i>	<u> 1</u>	REASON FOR CHANGE	OF CUSTOD	
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CHAIN OF CUSTODY RECORD ENVIRONMENTAL PROTECTION AGENCY REGION VII

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CONTENTS OF SHIP	MENT												
SAMPLE		<u> </u>	OF CONTAIN	NERS	VOA SET	+=			VEDIA oth	-	PECEIVING LABORAT	MATION	
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-303	3.	2,			/,	V	1]						
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DESCRIPTION OF S	HIPMENT				MODE OF S	HIPN	MENT	· 					-
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RELINQUISHED BY			E TIA	AE F	ECEIVED BY	_		`	/111-	7/	REASON FOR CHANGE	OF CUSTOD	Y
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RELINQUISHED BY	•	DAT	FE 111	ME	RECEIVED BY						REASON FOR CHANGE	. 0- 05101	, •
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HANG HE GUSTODY RECORD INTROHIBENTAL PROTECTION AGENCY REGION VII

ACTIVITY LEADER(PI	(Int)	Locurde	ey er activity (מושל ביי באור אינים או	su c	DATE OF COLLECTION 9/	SHEET 2
CONTENTS OF SHIPM	MENT			· • · • ·	DAY WONTH YEAR	10 11%
SAMPLE				SAMPLED MEDIA	RECEIVING LABORA	
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DESCRIPTION OF S	HIPMENT		MODE OF SH	IPMENT		
PIECE:S) C	ONSISTING			ERCIAL CARRIER _		
			CAMPL	ER CONVEYED	SHIPPING DOCUMENT	NUMBER)
PERSONNEL CUSTO			SCEIVED BY	. 34114	REASON FOR CHANGE	OF CUSTODY
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RELINQUISHED BY			CAR OBY		REASON FOR CHANG	OF CUSTODY
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RELINQUISHED BY	INSEALLE		TECELLED BY		REASON FOR CHANG	E OF CUSTODY
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HAIR OF CUSTODY RECORD SEVIRONMENTAL PROTECTION AGENCY REGION VII

ACTIVITY LEADER(PI		NAME OF SURVE		1 GK	35		1		COLLECTION		SHEE	
CONTENTS OF SHIPM		· 4-176.6 =1,E		-17.				DAY	MONTH YE		/ 100	<u>_</u>
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SAMPLE NUMBER	CUBITAINER		, CA SET , A S EA	1		dust		1	REMARKS OTHER	R INFORMA	ATION receipt.	
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			SAMPL	LER Ci	ONVL	EYED		ŝН	IPPING DOCUM	MENT N	UMBER)	_
PERSONNEL CUSTO	ODY RECORD	F FIME DE	ECEIVED BY			7,	,,-7	PEAC	ON FOR CHA	NGF C	FCLICTO	<u> کر</u>
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APPENDIX F MDNR WELL LOGS

Reported to Reserve ASLU	ľ
GEOLOGICAL SURVEY AND VATER RECOURCES	_
LOG NOD OWNER	
COUNTY FARM WELL NO	5
Sl. Louis Missouri Hills 2-8	
T B DRILLER	
4/7 76 DATE-Surd 15-Mont 68 Comp. 19.8-6	7
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800 5,00 OJ 1250 1300 osp.

APPENDIX G

ON-SITE DRILLING AND SAMPLING LOGS MARCH 4 THROUGH 9, 1991

P

ecology and environment, inc.

OVERLAND PARK, KANSAS

			clede Coal		Well Casing Size, Type, Quantity	/
Ste Lo	cation: S	t. Louis	, Missour	<u> </u>	Boring Diameter: 5 till	
Boring	Well Locati	on: Bot	1		Screen Size & Type: 1/4	
Boring/	(Well No.:	Bdi		(1-1	2) Screened Interval: 1/4	
Drilling	Contractor	John	Mathes &	Assoc.	Well Diameter:	
Dritting	Equipment	CME 55	O ATV S	plied Stent	Anger Date: Musch 4 1991	
Orlier:	Keith	Bunselm	eyer		Start Time: 13:05 Completion Time:	11:00
					Total Depth of Hole:Groundwate	
	_				Completion Depth:Surface E	
1-1 Ospin	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks
1,0					Gravel and Sund	HNU=NAB*
	BUI-A				Black gray clay with	
7.0	B\$1-8				Bluck lowing soil with gravel + sparre, lay bulls	NAB
7.0 -	-				Black loung soil with gravel and brown clay wells	NATS
-	- BOI-C				black gray clay	
75					medium y-riel own brick fragments	Tar-like odor
, 식 -					Block gray clay spaine angular gravel and brick	Tar-like odor Coal-ter like Odor only shee NAB
"	_	<i>;</i>			- Re En su! & ~ 421	
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OVERLAND PARK, KANSAS

emarks
= 2pm
= <u> </u>
= 20 mm
/ .^
= 60 pp.m bire hole



OVERLAND PARK, KANSAS

. 1	Project No	me & N	umper Lac	lede Coal	Gas	Well Casing Size, Type, Quanty	1/i+			
;	Borting/Well No.: 343					Boring Diameter: 5 / 1/2/				
						Screen Size & Type:				
						Screened Interval:// A				
						Well Diameter:				
	•		Bunselme			Start Time: 15:50 Completion Time:				
	Geologist	Wesl	ey McCal	1						
	Orller's H	elper(s):_	Jim Bar	ker, Jeff		Completion Depth:Surface Elevation:				
<i>‡</i> +,	O@DIT S	ampie	N-Value Number of Diows/6°	Gr aphic Log	Completion Date	Description	Remarks			
	-					Gravel and gray black				
2.0						Black chay loan soil with coal cinders	Heavy tur-like			
∄ ,Û							HNII = 2.5 jpm.			
.,,0						orth minor silt				
	-						HWH = Zppna			
3 15						Green pointie or ex	11N1 = 2 ppm			
: 6	_					Black only, tarry coze with coal cinders Gray Black Clay	1/114 - 1. ppm			
14	<u>-</u>					Gray Black Class	MAIN Ziponi			
11			,							
						Cemusal 2- 22				
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OVERLAND PARK, KANSAS

Project	Name &	Number: Lac	lede Coal	Gas	Well Casing Size, Type, Quanty 1/4	Well Casing Size. Type. Quanty		
Ste Loc	ation:S	t. Louis,	Missouri	<u> </u>	Boring Diameter: 5 : (Boring Diameter: 5 in (
			<u> </u>		Screen Size & Type: 1/4	Screen Size & Type: 1/H		
3oring/\	Well No.:_	23-			Screened Interval:			
			lathes & A	Assoc.	Well Diameter:			
Drilling	Ear Homeon	- CME 550	DATV 5	lid stein	1 1 1 19 Date: 11/1 1/2 14 1991			
			eyer					
		ley McCal				Start Time: 14:42. Completion Time: 21:54		
_				Crank				
Driller's	Helper(s):				Completion Depth:Surface Eleva	non:		
OPPIN	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks		
_					Grave (und see serry			
_					⊕9 ± (
					Gravel and bills			
					Grand and bick			
-	•				Fragments in gory sand	•		
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\ -	-				·			
-	-				Gravel with brick and wood fragments			
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ecology and environment, inc.

FT1307 FM00579SA

OVERLAND PARK. KANSAS

, ,	roject Nome & Number: Laclede Coal Gas					Well Casing Size. Type. Quanty				
9	ate Loca	mon: S	t. Louis	, Missouri	<u> </u>	Boring Diameter: 5 zuch	Boring Diameter: 5 Zuch			
	Boring W	/ell Locati	on: Byt.	5		Screen Size & Type:	NA			
i	Boring/W	el No.:	B. J. 5	·			NA			
					Assoc.	Well Diameter:	NA			
						11195 Date: 1116 4 1991				
			Bunselm	,						
	Geologis	: Wesl	ey McCal	1						
	Orllier's I	Heiper(s):_	Jim Bar			Completion Depth:Surface El				
<i>†</i> .	Ospin	Sample	N-Value Number of blows/6"	Graphic Lag	Completion Date	Description	Remarks			
. •	-					Sund and gravel with				
3,9	_				ž	gray bluib clay	slight coal-tur			
.ĉ	_ 									
	 					givery black clay	14 N'll = 0.4 ppm			
3.6	_					greenish quei clay.	HALL = C. + gp.m			
4.0			<i>;</i>			same as above				
23,							+NU = 10ppm			
	_					२०११र २२ ७०%				
10.1						bediock = limestone	Halil = Ippon down spent suger hole			



OVERLAND PARK, KANSAS

Project Name & f	Number Lac	lede Coal	Gas		NA		
Ste Location: S	t. Louis,	Missouri	·	Boring Diameter: 5 i.c./L	- Boring Diameter: 5 / 100/L		
Boring Well Locati	on: $B\phi$			Screen Size & Type:	NA		
Boring/Well No.:	BAL			Screened Interval:	NA		
Drilling Contractor	John M	lathes & A	ssoc.	Well Diameter:	114		
				inser Date: Marile 5 1991			
Orler: Keith	Bunselme	eyer		Start Time: 13/25 Completion Time: 13	:53		
Geologist: Wesl	ev McCall	1		Total Depth of Hole:Groundwate			
Driller's Helper(s):	Jim Barl	ker, Jeff	Crank	Completion Depth: $\underline{/\mathcal{V}}$ Surface Ele	evation:		
Osoli Sample	N-Value Number of biows/6*	Graphic Log	Completion Date	Description	Remarks		
			·	gravel and gray sandy			
				black lassy clay coil with gravel bizzk terry ooze	Hail > 5 pm		
				black tarry ooze oily sheen, brown clay so some some oravel	HNI ce Ipjime.		
				त्रह इंट्यार छात्रात राजार यज्ञाराष्ट्री	HN il = Spor		
	*			Augers broke Lotwer 2nd anger, unable to retrieve			

 $\Delta = \omega_{\rm eff} +$

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OVERLAND PARK, KANSAS

	Project :	Name & N	lumber. Lac	lede Coal	Gas	Well Casing Size, Type, Quanty	:		
						Boring Diameter: 5 inch			
	Borting V	Vell Locatio	on: Rdia	B		Screen Size & Type:	NH		
			RUL B			Screened Interval:	Nit.		
	Drilling (Contractor	John	Mathes & A		Well Diameter:	NA		
	Drilling !	Equipment:	CME 55	O ATV &	lid stem	Auger Date: Musch 5 1991			
			Bunselm	,		Start Time: 1/25 Completion Time: 1/2	1:24		
	Geolog	s: Wesl	ey McCal			Total Depth of Hole: 33 Grounawate			
	Driller's	Helper(s):_	Jim Bar	ker, Jeff	Crank	Completion Depth:Surface E	Completion Depth:		
i +.	Osolu	Sample	N-Value Number of piaws/6°	Graphic Log	Completion Date	Description	Remarks		
	_					Gravel, saul, brick			
5,0	=	. 				Green gray clay	wal two onlow		
5.0	_					·			
	-					Green gray clay with could and brick fragment	. · · · · · · · · · · · · · · · · · · ·		
16.6							HALL = 10ppel		
,,,,,	\ _					as above			
	_								
	;\ <u> </u>						Hill = 10, m		
			ŕ			plantie, with fragman	+3		
	-	- - -					dala se si		
5.0	: _	- <u> </u>			-		HNII = 12 pm		
	-	-				Greening area = 124			
کا رہ	· -	-				Breanish green clay			
	.	-					(matigue 1)		

'DRILLING & SAMPLING LOG



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CONTINUATION SHEET

roject Name & Number: Laclede Coal Gas	Date:	Morch	5	1441	

Ste Location: St. Louis, Missouri

Boring Well Location: $B \stackrel{\circ}{\to} B$

Oebin	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks
-					Compact greenish brown clay with trace of fine send - silt	
-					Brown stay saturated with black tarry sils	HAIII - Japan
					Refusal al 35 ft	
-				-	Bedicek !	
_						
-						
-						
_	-					

P

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OVERLAND PARK, KANSAS

roject Name & N	tumber. Lac	lede Coal	. Gas	Well Casing Sze, Type, Quanty	A/A
e Location: S	t. Louis,	Missouri	·	Soring Diameter: 5 inch	
ring Well Location	on: Rd	7		Screen Size & Type:	NA
ring/Well No.:				Screened Interval:	NA
		<u> lathes & A</u>	Assoc.	Well Diameter:	NA.
lling Equipment:	CME 550	O ATV So.	2: 8 stem	auger Date: Much 5, 1991	
Wer: Keith	Bunselmo	eyer		Start Time: 12:15 Completion Time:	
oologs: Wesl	ey McCal	1	·····	Total Depth of Hole: 29^{24} Groundwater	Depth:
riller's Helper(s):_	Jim Bar	ker, Jeff	Crank	Completion Deptin: Surface Ele	evation:
Samps Too	N-Value Number of blows/6"	Graphic Log	Completion Date	Description	Remarks
-				Gravel used brick.	
				framents in diger	HN'IL = N'AB
-				Black coze, coal-coke	
				fragments, wood fibers, greenish gray clay balls	
-					H 1/4 = 14 gpm
				Black : 02 c . v. th coke	79
				glad send wood tragments	
_				Black 102e with coke glad wind fragments space gray-green clay balls	40/12 10 .
				Bize 2 33=2 with coke +	71
			 	1 3 124 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	HNil = 7ppm
				dry, più die green clay	·
				Green clay with	
	1			Grain clay with	
-		1			ŀ
				Green gray clay so, the	
_		 		frailutá	`
	·			Crefusal at 29 ft on Line	stone bedrock

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OVERLAND PARK, KANSAS

roject Name & Number Laclede Co	oal Gas		anty	NA
ite Locamon: St. Louis, Missou	ıri	Boring Diameter:	ch	
Boring Well Location: B & 8		Screen Size & Type:		NA
Boring/Well No.: 668		Screened interval:		NA
Orilling Confractor: John Mathes	& Asapc.	Well Diameter:		NA
Orilling Equipment: CME 550 ATV	Solid stan in	ye Date: March 5	1991	
Oriller: Keith Bunselmeyer	•	Start Time: 15:20 Comple		;00
Geologs: Wesley McCall		Total Depth of Hole:	Groundwater	Depth:
Driller's Helper(s): Jim Barker, Je	ff Crank	Completion Depth:	Surface Ele	vation:
Sample N-Value Graphic blows/6' Log	Completion Date	Description		Remarks
		Abort L Shallow refi several p Brick + Co	riskl at	į.

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OVERLAND PARK, KANSAS

roject Name & No	umber: Lac	lede Coal	Gas		1/4
te Location: St	Louis,	Missouri		8oring Diameter: 5 :::::	
oring Well Locatio	n: 13/1			Screen Size & Type:	N/A
oring/Well No.:	1399			Screened Interval:	NA
		lathes & A	ssoc.	Well Diameter:	1/ A
xilling Equipment:.	CME 550	ATV, 😙 .	1 sten a	March 5, 1991	
muer: Keith	Bunselme	yer		Start Time: 16.08 Completion Time:	4155
Geologist:_Wesle	ev McCall	<u> </u>		Total Depth of Hole:Groundwate	Depth:
Oriller's Helper(s):	Jim Bark	ker, Jeff	Crank	Completion Depth:Surface Ele	evation:
Osola Sample	N-Value Number of Diows/6°	Graphic Log	Completion Date	Description	Remarks
				Abort Location, Shallow refusal at	~
				Shallow refusal at	
				several points, Concrete foundation	
-				Concrete foundation	
				•	
-					
	<i>;</i>				
				,	
-					
, ,	1	1	i	1	1



OVERLAND PARK, KANSAS

roject Name & Number: La	clede Coal	Gas		anny UH	
te Location: St. Louis	, Missouri		Boring Diameter:	<i>:</i> 1	
oring Well Location: $\frac{\beta}{1}$	<i>‡</i>		Screen Size & Type:	NA	
oring/Well No.: B/#			Screened Interval:	NA	·
rilling Confractor: John	Mathes & A	ssoc.	Well Diamy ter:	N.A	
orilling Equipment: CME 5	50 ATV 5:/	a stea no	ysts Date: Merch 5	1991	
Miler: Keith Bunseli	neyer		Start Time: 7:05 Comple	otion Time:	35
Geologist: Wesley McCa	11		Total Depth of Hole:	Groundwater Dep	om:
Critier's Helper(s): Jim Ba	rker, Jeff	Crank	Completion Depth:	Surface Elevation	n :
Osó _{li} Sample Namber of NAdine	Grapnic Log	Completion Date	Description		Remarks
-			Absit Location		
			shallow refusal	at	
			several points.	,	
			Concrete Found		
_					
_					
			•		
				1.	

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OVERLAND PARK. KANSAS

-	Project Name &				Well Casing Size. Type. Quanty	•			
					Boring Diameter 5 inch				
	Boring Well Loca		<u>!</u>		Screen Size & Type:	Screen Size & Type: VA			
	Boring/Well No.:.	RIL		- 	Screened interval:/	L'A			
	Drilling Contract	or: <u>John</u>		_	Well Diameter:/				
	Drilling Equipmen	nt: CME 55	D ATV 50/	d stem	ruger Date: 11/20-16 6 19-91	<i>!</i>			
	Driller: Keit				Start Time: <u>05:10</u> Completion Time:				
	Geologist: Wes				Total Depth of Hole:				
	Driller's Helper(s)				Completion Depth: Surface El				
24.	Oppir Sample	N-Value Number of blows/6*	Graphic Log	Completion Date	Description	Remarks			
	_				Gowith and siend,				
<u>,</u>	-			 	Medium brown lay wil				
					medium brown = lay				
5					, , , , , , , , , , , , , , , , , , , ,				
	-				medium brown clay with sparse brick fragments	-			
					partie brick tragments				
1;				 					
					brick frayments				
13					gray black clary, wet				
٠ ٢					Reduced x 1524	HIV'U = 85 pm down open bering			
		·							
	-								
	-								
		1							



OVERLAND PARK, KANSAS

roject Name &	Number: Lac	lede Coal	Gas	Well Casing Size, Type, Quantity	NA
te Location:	St. Louis	Missouri		Boring Diameter: 5 Lick	
				Screen Size & Type:	
				Screened Interval:	
				Well Digmeter:	
rillina Eaulomer	CME 550	O ATV Sol	d stem .	ager Date: March 6 14	9/
xiler:Keit				Start Time: 07:05 Completion Time:	
				Total Depth of Hale:Grounawate	
				Completion Depth:Surface E	
Osolia Samble	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks
-	-			Gravel and sund	
-				Brown clay loam	
				Greatish arm clay	HALL : NAB
				Greatic greenish gray clay, damp, vily sheen	
				Greenish gray clay with	coal to-
				Refusal at 11 H	HNU > 50 ggm down spell bore hole.
	<i>;</i>				

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OVERLAND PARK. KANSAS

-	Project Name & I	Vumber La	clede Coal	Gas	Well Casing Size. Type, Quanty	NA			
	Ste Location: S	t. Louis	, Missouri	<u> </u>	Soring Diameter: 5 inch				
	Boring Well Locati	on: <u>B13</u>	- ,		Screen Stze & Type:NA				
	3aring/Well No.:_	B13			Screened Interval:	N' A			
	Drilling Contractor	John	Mathes & A	Assoc.	Well Diameter:	NA			
					anger Date: March 4				
			,		Start Time: <u>09: 45</u> Completion Time:				
					Total Depth of Hole:				
					Completion Depth:Surface				
ÿ. H.		N-Value Number of Diows/6*	Grapnic Log	Completion Date	Description	Remarks			
ŧ					Errel ind Ford				
10					Poor recovery - Wet black 2024 of h whendon't brick fragmen black wet soze brick fragments Foor Recovery Gray Black clay, brick fragments black coze	-			
		,			Refusel n+ 12 ft	Haining, no			

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OVERLAND PARK, KANSAS

Project Name	& Number Lac	lede Coa	l Gas	Well Casing Size, Type, Quanty	<u> </u>
Ste Location:	St. Louis,	Missour	<u>i</u>	Boring Diameter: 5 Zuch	
				Screen Size & Type:	N/A
Boring/Well N	o: 1314	 		Screened Interval:	NA
Drilling Contro	scror: John l	Mathes &	Assoc.	Well Diameter:	NA
Drilling Equipr	nent: CME 550	O ATV Sil	id stem a	1945 Date: March 6, 1991	
		,		Start Time: 12:35 Completion Time: 1/	•
Geologist:\	eslev McCal	1		Total Depth of Hole:Groundwate	er Depth:
	or(s): Jim Bar			Completion Depth:Surface E	•
Oppir Sam	N-Value Number of blows/6*	Grapnic Log	Completion Date	Description	Remarks
				Gravel and sandy soil	
-				livy little receiving.	
				Black oil on surer	
 				with traces of soil,	
				gravel and brick	
			- 	Pragman 15	•
				,	
				Pery little receivered	
-				over entire internal,	MALL : 40 pp.
				•	
				Blade vil or sugars.	
				Sparse clay and brick	Sub sur Pace
-				Freignettic 1 to	rusole may
				\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	he stripping
-					ing at si
		1		1 1	Screens ty stemp e
-				Cerusat +	I of oil collected
					ut = 3 ++ dept

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OVERLAND PARK, KANSAS

P	roject Name	& N	umber: Lac	lede Coal	Gas	Well Casing Size, Type, Guanty	NA
S	ite Location:	St	Louis,	Missouri	<u> </u>	Boring Diameter: 5 Lich	
9	loring Well Li	ocatic	n:	5		Screen Size & Type:	NA
=	loring/Well N	o.:	B15			Screened Interval:	1/t
1	Oriiling Contro	actor:	John_N	lathes & A	Assoc.	Well Diameter:	NA
ſ	Ortilling Equipm	nent:	CME 550	ATV (lid stem	anger Date: Much 6, 1991	
Į	Oriller: Ke	ith	Bunselme	yer		Start Time: 14:05 Completion Time: 10	1:50
			ey McCal			Total Depth of Hole: 31 ff Groundwater	
1	Oriller's Helpe	er(2):	Jim Barl	ker, Jeff	Crank	Completion Depth:	vation:
1	OBOIN Sam	pie	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks
	_	-				Gravely brown clay loan soil Greensh crom city or the Black term in a terral	
	-		Ì			Breenish krom ody 15th	
•						Black tar sutunted soil	
	_				·	Black for saturated	
						Black for saturated sand, some greensh	•
t						1) ey clay 5x/16	
-							
	_					•	
5				<u></u>			
Ø						Black for saturated sand, brick fragments	
						o, ly avent	HNU = 20,000
O			i i				
<u>.</u>						Bizzk terry sew (ust(1)	_
5						Gray black clay	
	-					Black, wet word Evers	
.'7						brick fragments.	we+

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OVERLAND PARK, KANSAS

roject Name & Number: La	clede Coal	Gas	Well Casing Size, Type, Quanty	1/A
te Locamon: St. Louis	, Missouri		Boring Diameter: 5 7/1//	
oring Well Location: $3/$			Screen Size & Type:	!/A
oring/Well No.: <u>Bib</u>			Screened Interval:	NA
Contractor: John			•	NA
Orilling Equipment: CME 55	50 ATV , 52/3	1 steel a	Wer Date: 11/a-ch 6, 1991	
Oriller: Keith Bunsel	neyer		Start Time: 15/05 Completion Time:	5:24
Secogs: Wesley McCa	11		Total Depth of Hole:Groundwate	r Depth:
Oriller's Helper(s): Jim Bar	rker, Jeff	Crank	Completion Depth:Surface El	evation:
Oso Sample Number of Diows/6"	Graphic Log	Completion Date	Description	Remarks
			Repealed refusal at a 3 Rt depth on foundation. Abort Location	HNU - 150 gp. a Noun 3 Kt hole

P

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Project	Name &	Number: La	clede Coal	l Gas	Well Casing Sze. Type. Quanty	NA		
					Boring Diameter: 5 in all			
					Screen Size & Type:	· · · · · · · · · · · · · · · · · · ·		
		B17				N/A		
						# / . \		
					Well Diameter: inger Date: Micric 6, 1991	_//		
			•			/ / / 0		
					Start Time: 15:138 Completion Time: 1			
					Total Depth of Hole: $\frac{35}{9}$ Groundwate			
Driller's	Helper(s):	Jim Bar	ker, Jeff	Crank	Completion Depth:Surface Ele	evation:		
OSOIL	Sample	N-Value Number of biows/6*	Grapnic Log	Completion Date	Description	Remarks		
					Black sunder pochles Black learn, soil with coal + cinders			
[] =	•				Black learn, soil with coul + cinders			
. =					Gray black clay, sparine brick fragments and cont distants			
· -					Sundy you black clay			
-	-				with sparse horice	<u>-</u>		
. _					Sundy gray black clay with sparse brick fragments	HNU = NAB		
	-				Sanda graevinh black class			
-	- -				Brounish clay with abundant			
s	-			 	angular gravel			
-	- - -				Gray black clay onth			
_ _	_							
-	-				Sundy green black clay	HNU = 4 ppm		
_ -					pabbles	plenotex = Oggm		
` -					Sandy freen slack clay			
	_				proportion of assembly	HN'll = 35 pm		

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OVERLAND PARK, KANSAS

P	roject !	Name & M	Number Lac	lede Coal	Gas		- Well Casing Size. Type. Quanty			
Ste Locamon: St. Louis, Missouri						Boring Diameter: 5 inch				
2.1						Screen Size & Type:	NA			
É	oring/V	Veli No.:	BIS			Screened Interval:	1 t			
			John M	lathes & A	Assoc	Well Diameter:	11,+			
í	Oriting (Equipment	CME 550	ATV 50 C	id Stem	Auger Date: March 6 1991				
			Bunselme	,		Start Time: 16:50 Completion Time: 17	/ 			
(Geologi	s: Wesl	ev McCall	<u> </u>		Total Depth of Hole: 33 4 Groundwater				
į	Orllier's	Helper(s):_	Jim Barl	ker, Jeff		Completion Depth:Surface Ele	,			
	Ospin	Sample	N-Value Number of blows/6"	Graphic Log	Completion Date	Description	Remarks			
	_					Brown sandy soil with				
į	_			·····		Blook sandy clay, parte angular gravel	HNU = 110 ppm			
						Black clay, course soud and sparse ungular gravel	-			
^	_					Wet gray black clay with	HN11 = 500 ppm			
•						coarse to medium sand				
Ĺ							Hh.d = 200 ggmi			
12	-		·			Breenish black clay				
	-	•								
	-	-					14 Nil = 130 gpin			
5						Greenish black clay	= 14/3 (= 1/4/3			
	1.	_ 1	1	I	ĺ	1	t			



OVERLAND PARK, KANSAS

Project Name & Number: Lacle	ede Coal Gas	Well Casing Size. Type. Quanty/	- Well Casing Size. Type. Quanty		
Ste Locamon: St. Louis, N	Missouri	Boring Diameter: 5 / / /			
Boring Well Location: 13]	1	Screen Size & Type:	N'A		
Boring/Well No.: B19		Screened Interval:	1.'A		
Drilling Contractor: John Mat	ì		n'A		
Ortiling Equipment: CME 550	ATV, sold stem a	1985 Date: March 1, 1991			
Oriller: Keith Bunselmey		Start Time: <u>194:20</u> Completion Time:7			
Geologs: Wesley McCall		Total Depth of Hole: $\frac{30}{20}$ Groundwater	Depth:		
Oriller's Helper(s):Jim_Barke	r, Jeff Crank	Completion Depth:Surface Ele	votion:		
Osolo Sample Number of Diows/6*	Graphic Completion Date	Description	Remarks		
-		brown sundy lound soil agreed			
		Black sludge of coal conders, stay, and spaine horized. Fray ments	14 Alil = 14 M		
		Black sandy clay with timestone gravel and could conder so slay Det black sand?) abundant roal clinders and borick fragments.	MNU - NAB		

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Project Name & Number: Laclede Coal Gas	Well Casing Size, Type, Quanty
Ste Locamon: St. Louis, Missouri	Boring Diameter: O Charle
Boring Well Location: $BZ\phi$. Screen Size & Type:
Boring/Well No.: BZ#	Screened Interval:
	Well Diameter:
Drilling Equipment: CME 550 ATV solid stem auger	Date: March 17, 1991
Orller: Keith Bunselmeyer	Start Time: <u>09:19</u> Completion Time: <u>10:20</u>
Geologst: Wesley McCall	Total Depth of Hole: 30PF Groundwater Depth:
Orller's Helper(s): Jim Barker, Jeff Crank	Completion Deptin:Surface Elevation:

Sam	N-Value Number of Number of	Graphic Log	Completion Date	Description	Remarks
				Coal waste.	
				Sandy coul cinders and blog, some orisk	
				blay, some orisk	
				frag wents and gravel	
_					HN'Il = NA B
					7
_					
		<u> </u>		us above with wood fragments + fibers	
	<i>,</i>			as above: Wet	
-				Sundy gray black clay	

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OVERLAND PARK. KANSAS

Project Name & Number Laclede Coal Gas	Well Casing Size, Type, Quanty
Ste Locamon: St. Louis, Missouri	Boring Diameter: 5 inch
Boring Well Location: 121	Screen Size & Type:
Boring/Well No.:	Screened Interval:
Oriling Confractor: John Mathes & Assoc.	Well Diameter:
Drilling Equipment: CME 550 ATV	Date: 1100ch 17, 1991
	Start Time: 10:30 Completion Time: 4:40 11:10
Geologist: Wesley McCall	Total Depth of Hole:Groundwater Depth:
Orller's Helper(s): Jim Barker, Jeff Crank	Completion Depth:Surface Elevation:
Sample N-Value Stappic Complete Dots	
	shallow refinal (43Ht) at zeveral points. Limestone hadrock. About Location

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oring Well Location: 3 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	Missouri Mathes & As O ATV	SBOC.			
Oriller's Helper(s): N-Value Number of blows/6*	Graphic Log	Completion Date	Completion Depth:Surface Description	Remarks	
			it swa-al points. Line: Jone astrock. About Licution.		

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OVERLAND PARK, KANSAS

ject Name & Number: Lacledo	e Coal Gas		onty N	<u>/\</u>
Location: St. Louis, Mi	ssouri	Boring Diameter: inc	<u>k</u>	
ing Well Location: 1323		Screen Stze & Type:		
ing/Well No.: B23		Screened Interval:		A
ing Contractor: John Math				A
ing Equipment: CME 550 AT				
	,	,	· ·	
er: Keith Bunselmeyer		Start Time: //:-*** Comple		
Ologist: Wesley McCall		Total Depth of Hole: $\underline{\hspace{1cm}}^{\mathcal{Y}}$		
ler's Helper(s): Jim Barker,	Jeff Crank	Completion Depth:	Surface Elevation:_	
	consistion cog Date	Description		Remarks
		2200 (0000 400) on		
	Ke	d brown sandy co	lay leave	
		enthered bedrock	[· · · · · · · · · · · · · · · · · · ·
-		<u> </u>		
		Refusal a-	1 24 FA	-
		Limestone	hedreck	
-		• • • • • •		-
-				
		•		
- ;				
_				
-				
			1	

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OVERLAND PARK, KANSAS

	Project i	Name & 1	Number Lac	clede Coal	. Gas	Well Casing Size, Type, Quanty	NA			
	Site Loc	ата п :S	t. Louis	, Missouri	· 	Boring Diameter: 5 /nc/	- Boring Diameter: 5 /nc/c			
	Boring V	Veti Locat	ion: B2	4		Screen Size & Type:	N'A			
	Boring/V	Vell No.:_	B24		 .	Screened Interval:	NA			
	•		John J	Mathes & A	Assoc.	Well Diameter:	NA			
	Drilling (Equipment	CME 55	0 ATV 50 (id stem a	2481 Date: 11 sich 7, 1991				
	_		Bunselm			Start Time: 1313 C Completion Time: _/*	4:05			
	-					Total Depth of Hole:	er Depth:			
	Driller's	Helper(s):	Jim Bar	ker, Jeff	Crank	Completion Depth:	Hevation:			
1	Osoin	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks			
1	_					Gravel , brown sundy sil				
						Grassl brown sandy with bragish green clay with sparse coal-coke trogments				
5					-	Grayish green clay, damp, plantic				
							HNU = NAB			
1 %	._									
i E						-traces of coal day	coal law odos HNil = NAB			
2		-				Sandy year way sing	/			
2 2	2 -	-	<u> </u>			Sandy item in a sup sing the frage frage for super frage mater all with court courts	slight coal top alor			
2.						Refusal not 23ft	7.77			
	1	ì	1	1	ì	1	• 1			

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OVERLAND PARK. KANSAS

	Project (Name &	Number: La	clede Coal	l Gas	Well Casing Size, Type, Quanty	11:4		
						8oring Diameter: 3.75 ID	_		
	Boring V	Vell Locat	10n: <u>B2</u>	′/		Screen Size & Type:	7/4		
			355			Screened Interval:	NA		
				Mathes & A	Assoc.		2 A		
	Drilling (quipmen	. CME 55	0 ATV, 401	Vew stem	Hayer Date: 11/25-2 17, 17-11			
			n Bunselm			Start Time: 14:117 Completion Time: 17	125		
	Geologi	s: Wes	lev McCal	1		Total Depth of Hale: $\frac{2^{3}}{2}$ Groundwate	r Depth:		
	Driller's	Helper(s):,	Jim Bar	ker, Jeff	Crank	Completion Depth:Surface El	evation:		
<i>!</i> .	Ospin	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks		
1	-	×				Sand and Grave			
3		55*				Sand and Grave. Green brown randy clay with gravel + sparse and Frayments			
-	_	55				No Recovery	No Recovery		
5	-	55		_		Brown + black mottled ciay spurse coal + gravel fragments Brown Clay			
7	+=	4-				Gray Islack Clay with coal cind	a''4		
9		55		 -		Black mottled green clay	-10% Recovery		
1	_					Greenish cray lay	-40% Recovery		
	=	55				with black nottling (Manganese)	-15% Recovery		
	-						/) il		
1.11	=			,			-30% Decelery		
14	 	55	;			Compact your clay, 49thy	Layon Receivery		
21		5.5				black motther green chap	1 40% Receivery		
3						Line grained crystalline lime stone = verticib			

FT1307 FM00579SA

OVERLAND PARK, KANSAS

Project N	vame & I	Number: Lac	lede Coal	Gas	Well Casing Sze. Type, Quantity	is 4
Ste Loco	ation: S	t. Louis,	Missouri	L	Soring Diameter: 4,25 fact Z	ת
Boring W	/ell Locati	on: B&	b		Screen Size & Type:	NA
3oring/W	/eli No.;	C32			Screened Interval:	1/4
			lathes & A		Well Diameter:	MA
					1163 AT Date: March 1/ 1991	
		Bunselme	,		Start Time: <u>17:01</u> Completion Time: <u>10</u>	:12
Geologs	w: Wesl	ey McCal			Total Depth of Hole:	
_			ker, Jeff		Completion Depth:Surface El	
Ospin	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	Remarks
	_				Gravel , took to table.	No sumple
	<i>(</i>				Black sandy studge with city sheen some gravel fragments	(n 20%) HNU = NHB
	<u>C</u> S				Green compact clay sparse areas of las condamination	(245%)
	CS				Green compact clay with black mangarage	(پوجو و سد)
		,			Half borney chainsensen	HNU = 25 ppm down open sore hole

4

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OVERLAND PARK, KANSAS

P	roject i	Name & I	Number:_Lac	clede Coal	Gas	Well Casing Size, Type, Quanty	.21A
Ste Location: St. Louis, Missouri						Boring Diameter: - 25 /n:/	エ カ
В	oring V	Vell Locati	on: Bit	4		Screen Size & Type:	4/4
2 2 2						Screened interval:	VA
				Mathes & A		Well Diameter:	NA
						auger Date: 1/15/ 8, 1991	
٤	oriler: _	Keith	Bunselm	eyer		Start Time: 113:115 Completion Time: 113	31/
(Seologi	st: Wesl	ey McCal	1		Total Depth of Hole:Groundwate	Depth:
ί	Criller's	Helper(s):	Jim Bar	ker, Jeff	Crank	Completion Depth:Surface Ele	evation:
	Osolk	Sample	N-Value Number of blows/6°	Grapnic Log	Completion Date	Description	Remarks (*c Recovery)
	_	_				Gravel, Fock in black spoil, taking 20005	
		25				•	of sampler.
i	_	55				Green compact clay with zones of tar contamination	, ,
			,			Black farry clay Green (empact Chay 14 it Halt noticy advancement	HAIL = 7 ppm NO contains had there ed
	=	-					•

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OVERLAND PARK, KANSAS

				L Gas	Well Casing Size, Type, Quanty	NA		
				<u>i</u>	Soring Diameter: 4.25 i.i.ch I	- Soring Diameter: 4,25 i.u.h I.D.		
orting V	$\circ u$				Screen Size & Type:			
oring/V	Vell No.:	654			Screened Interval:	1'A 1'A		
			lathes &	Assoc.	Well Diameter:			
		Bunselme			Start Time: 12:40 Completion Time: 16	- Date: March 13 1991		
		ey McCal			Total Depth of Hole:Groundwate			
_			ker, Jeff					
					Comptetion Depth:	levation:		
Ossil	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	(7. Remarks		
	-				Gravel, rubble, smuly exil	1/0 Swaple		
_	55				Black taily sandy clay soil brick + wood flag ments Sand, gould + crick fragments suturated with greenish black oil			
	35				Sand, gravel + crisis	(~10%)		
_	55				greenish black oil	(~26%)		
				 	Augers conted with oil	HAVU = 9 ppm		
	55					1		
	\ \				Brick fragments, gravel + sent	(45%)		
=	1				Brick and concrete fragments in sand, saturated with oil	(410 %)		
					Refusal at 14 ft.	HNU = 13 mm Lour spin hore		
-		<i>;</i>		·		Sample of oil and water collection applit apoint		
	- - -							

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OVERLAND PARK. KANSAS

Project Name & Number Laclede Coal Gas				Gas	Well Casing Size. Type. Quanty	1/4
Ste Location: St. Louis, Missouri					Boring Diameter: 4,25 inch	三の.
Boring/Well No.: C 5.55					Screen Size & Type:	NA
					Screened Interval:	1/A
					Well Diameter:	Λ'A
-						
	Keith	Bunselm	PVPT		Start Time:Completion Time:	111 + 3 (24)
						
_		Ley McCal				-
Oriller's	Helper(s):	JIM DAL	KEI, JEII	Olank	Completion Depth:Surface E	evation:
Ospin	Sample	N-Value Number of blows/6"	Graphic Log	Completion Date	Description	Remarks
_					Gracel and brick tragment in black soil	13 No Sample
	55				Mark Shark Combas.	19572) 41/4 = NA-B (~ 2090)
-	55				Dariz green samy 5/24	(n 20%)
=	5.5				Green black clay brick and limes ton frayments, lans it conting and many politics	(~50%) 4111 = 260 ppm
					Rifusal it split s	

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OVERLAND PARK, KANSAS

Project Name & Number: Laclede Coal Gas				. Gas		nA		
Ste Location: St. Louis, Missouri					Soring Diameter: 425 inch	1		
Baring Well Location: B \$\frac{1}{7}\$					Screen Size & Type:	11.4		
Boring	/Well No.:	C56			Screened Interval:	NA		
Drilling	Contractor	John J	Mathes & A		Well Diameter:	NA		
					Date: March 8. 1991			
		Bunselm						
Geol	ogs: Wes]	ley McCal	1					
Orition	's Helper(s):	Jim Bar	ker, Jeff	Crank	Completion Depth:Surface El			
O _O O	Sample	N-Value Number of blows/6°	Graphic Log	Completion Date	Description	(70 Recovery)		
-	_					No Sample: Bore hole		
-	1/5				•	Bire hole adjacent to		
-	-		٠.			CS5		
-	_							
-	<u>-</u> ss				course from and Francis	·		
	- 55				green clay botton (inches 4 ternation cavers of sund + gravel and green clay	cil suturated		
	55				(carse sand, medium rounted	(10%)		
. -	55				Wied fibers, black torry sands green compact clay last Zina	(10%)		
		;			Sald Aresement of			
	- - -					·		
-	_							

FT1307 FM00579SA

OVERLAND PARK, KANSAS

Project Name & Number Laclede Coal Gas				Gas	Well Casing Size, Type, Quanty	1/4		
Ste Locamon: St. Louis, Missouri					Boring Diameter: 125 1/2/ I	- Soring Diameter: 15.25 inch ID		
בו כו					Screen Size & Type:	1/A-		
Boring/	Well No.:_	<u>CS7</u>		· · · · · · · · · · · · · · · · · · ·	Screened Interval:	1/1		
Drilling	Contracto	. John 1	lathes & A	ssoc.	Well Diameter:	NA		
Drilling	Equipment	CME 550	DATY , 1/c/	low stem A	fuger Date: 1/25-1/9 1991	_ Date:		
Ortier:	Keith	Bunselme	eyer		Start Time: 03 + Completion Time: 09			
Geolo	gs: Wes	ley McCal	1		Total Depth of Hole: 12.5 Groundwate	Total Depth of Hole: 12.5 Groundwater Depth:		
Orlier's	Helper(s):	Jim Barl	ker, Jeff	Crank	Completion Depth:Surface Ele	evation:		
OSOIT	Sample	N-Value Number of blows/6"	Graphic Log	Completion Date	Description	Remarks		
	1/5					No Sample		
-	55			_	soudy green-black clay	(50%)		
-	55				ord brick	(40%)		
-	55				brick frequents	(10%) 1'hoto 3, coll 2		
-	-				Bisch wet with brick fragments + gravel Black Sund, Fine subjournment	Photo 9 roll 2		
5 =				 	gravel c	77.76-		
-	-				Refusal at 12.5 ft	Hail = 3pp.n down open bore hole		
-	- -					Sole Note		
-	_	,						
-	_	,						
-	-							
-	-							
.	-							
.	-							
'								